



PUBLISHED FOR SISSA BY SPRINGER

RECEIVED: September 23, 2013

REVISED: February 12, 2014

ACCEPTED: March 5, 2014

PUBLISHED: April 3, 2014

Dimension changing phase transitions in instanton crystals

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ABSTRACT: We investigate lattices of instantons and the dimension-changing transitions between them. Our ultimate goal is the $3D \rightarrow 4D$ transition, which is holographically dual to the phase transition between the baryonic and the quarkyonic phases of cold nuclear matter. However, in this paper (just as in [1]) we focus on lower dimensions — the 1D lattice of instantons in a harmonic potential $V \propto M_2^2 x_2^2 + M_3^2 x_3^2 + M_4^2 x_4^2$, and the zigzag-shaped lattice as a first stage of the $1D \rightarrow 2D$ transition. We prove that in the low- and moderate-density regimes, interactions between the instantons are dominated by two-body forces. This drastically simplifies finding the ground state of the instantons' orientations, so we made a numeric scan of the whole orientation space instead of assuming any particular ansatz.

We find that depending on the $M_2/M_3/M_4$ ratios, the ground state of instanton orientations can follow a wide variety of patterns. For the straight 1D lattices, we found orientations periodically running over elements of a \mathbb{Z}_2 , Klein, prismatic, or dihedral subgroup of the $SU(2)/\mathbb{Z}_2$, as well as irrational but link-periodic patterns. For the zigzag-shaped lattices, we detected 4 distinct orientation phases — the anti-ferromagnet, another abelian phase, and two non-abelian phases. Allowing the zigzag amplitude to vary as a function of increasing compression force, we obtained the phase diagrams for the straight and zigzag-shaped lattices in the (force, M_3/M_4), (chemical potential, M_3/M_4), and (density, M_3/M_4) planes. Some of the transitions between these phases are second-order while others are first-order.

Our techniques can be applied to other types of non-abelian crystals.

KEYWORDS: AdS-CFT Correspondence, Phase Diagram of QCD, Holography and condensed matter physics (AdS/CMT)

ARXIV EPRINT: [1304.7540](https://arxiv.org/abs/1304.7540)

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1 Introduction

In the ordinary $N_c = 3$ QCD cold nuclear matter forms a quantum liquid, but for large N_c it becomes a crystalline solid [2]. In many holographic models of QCD [3, 4] baryons are represented by instantons of the $U(N_f)$ gauge theory living on the flavor branes [5], so cold nuclear matter corresponds to a whole crystalline lattice of instantons [6]. The geometry and even the dimensionality of this lattice depend on the baryon density: at medium densities, the instantons form a 3D lattice in the $x_4 = 0$ hyperplane of the holographic 4D space (not counting the time); this corresponds to the baryonic phase of nuclear matter. At high densities, the baryons spread out into the x_4 dimension and form a 4D lattice; this is dual to the quark liquid phase of nuclear matter, or rather to the high-density *quarkyonic* phase [19] in which the quarks fill the Fermi sea, but the excitations near the Fermi surface are baryon-like bound states of N_c quarks rather than free quarks by themselves. According to Rozali et al. [7], the chemical potential of the quarks in this phase is related to the thickness of the lattice in the x_4 direction.

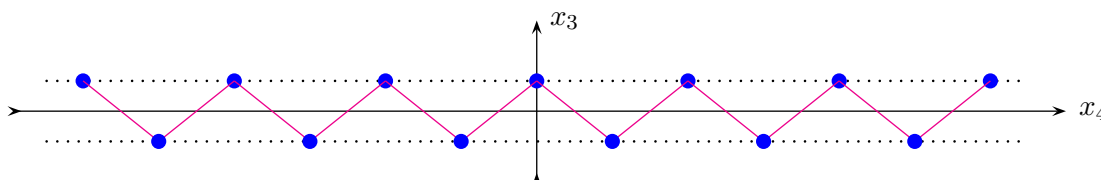
Most investigations of the low-temperature holographic nuclear matter focus on the extreme high-density limit where the x_4 dimension is thick and can be studied macroscopically (for example, see [7–15]). But our main interest is in the microscopic structures of the instanton lattices, and especially the phase transition between the 3D and 4D lattices. In our previous paper [1] (with Dmitry Melnikov) we found that there is a whole sequence

of such transitions: from a single 3D layer in the $x_4 = 0$ hyperplane, to two 3D layers in two parallel hyperplanes, to 3 layers, to 4 layers, etc., until the number of layers becomes too large to count individually. Alas, the 3D and 4D instanton lattices turned out to be too hard to analyze, so we resorted to oversimplified toy models. In the first toy model, we approximated the instantons as point-like charges repelling each other with 4D Coulomb forces; this approximation is very crude — it ignores the instantons' orientations, never mind the interference between the instantons — but it makes for a simple model of transitions between lattices of different dimensions. Indeed, in this model we found that increasing density makes a 3D lattice of instantons in $x_4 = 0$ hyperplane suddenly split into two sublattices in the $x_4 = \pm\epsilon$ hyperplanes; pictorially, this looked like the popcorn suddenly jumping up in the popper, so we dubbed the dimension-changing transitions the *popcorn transitions*.

In our second toy model, we used actual instantons and the exact ADHM [16] construction for the multi-instanton system. However, we had confined the instanton centers to a 2D plane $x_1 = x_2 = 0$ by making the inverse 5D gauge coupling rise steeply in the x_1 and x_2 directions. We also had the $1/g_5^2$ slowly rising in the x_3 direction, altogether

$$\frac{8\pi^2}{g_5^2(x)} = N_c \lambda M \left(1 + M^2 x_3^2 + M'^2 (x_1^2 + x_2^2) + O(M^4 x_{1,2,3}^4) \right) \quad \text{for } M' \gg M. \quad (1.1)$$

Consequently, at low densities the instantons formed a 1D lattice in the x_4 dimension (which acted as the model's only flat space dimension), but for higher densities the instantons moved into the x_3 dimension (which acted as the holographic dimension of the model). In that model, we saw two phase transitions in response to increasing instanton density. The first transition — from a straight 1D chain of instantons to a zigzag-shaped chain



— is second-order, so the zigzag amplitude behaves as $\epsilon \propto \sqrt{\rho - \rho_c}$. On both sides of this transition, the instantons' orientations form an anti-ferromagnetic pattern: two alternating orientations related by a 180° rotation. At higher densities there is another transition, but this time it's first-order: the zigzag amplitude ϵ jumps discontinuously, and the instantons' orientations also change their pattern: the orientations of the nearest neighbors are now related by a rotation through angle $\phi < 110^\circ$ instead of 180° .

Presumably, at still higher densities we would have seen popcorn transitions from the zigzag — which is a kind of two-layer lattice — to a 3-layer lattice, then to 4 layers, etc., etc. However, calculating the net energies of such multi-layer lattices turned out to be too hard, so we stopped at the zigzag.

Technically, the main difficulty in working with more complicated instanton lattices is setting up the ADHM construction [16] for the infinite number of instantons. Indeed, the ADHM construction for A instantons involves 4 non-commuting $A \times A$ matrices Γ_{mn}^μ whose

off-diagonal elements follow from non-trivial constraint equations; for $A \rightarrow \infty$ solving these equations becomes very hard. For a simple 1D lattice of instantons whose orientations are related by *commuting* $SU(2)$ symmetries, the ADHM constraints were solved by Kraan and van Baal [17], and we have re-derived and used their explicit formula for the instanton number density profile $I(x)$ in [1]. For a zigzag-shaped instanton chain with link-periodic instanton orientations, we found an exact solution of the ADHM constraints in terms of Fourier transforms of some rather messy functions, but we could not use it to derive an exact formula for the $I(x)$; instead, we had to use the small-instanton-size approximation. And it gets worse for more complicated instanton lattices and/or orientation patterns: for the *simple* 2D and 3D lattices and purely abelian orientation patterns the ADHM constraints have known formal solutions¹ — which alas are way too complicated for any practical use — while for the non-abelian orientation patterns there are no known solutions at all. This is particularly unfortunate since the lowest-energy 3D configuration for low densities is likely to be an FCC lattice with non-commuting orientations of the instantons [18].

In this paper we follow a shortcut around the ADHM construction — and also around the more difficult part of the energy calculation. We assume from the beginning that the instanton density is not too high, so the distances between instantons are much larger than the instantons' radii. In section 3 we show that for all such multi-instanton systems, interactions between the instantons are dominated by the two-body forces. The irreducibly-three-body forces, 4-body forces, etc. also exist, but they are suppressed by powers of the small (radius/distance)² ratio compared to the two-body forces. Consequently, we obtain a manageably simple approximate formula for the net energy of a multi-instanton system in terms of the instantons' positions and orientations.

Thanks to this formula, we do not need to assume (as we did in [1]) that in a lattice of instantons, their relative orientations have the same symmetries as their relative positions. Instead, we may use numerical methods to minimize the net energy with respect to all the orientations treated as independent parameters, without any symmetry assumptions. The algorithm we use in this paper is fairly simple: we set up a large lattice (200 sites is more than enough for 1 dimension) of fixed geometry, with a free $SU(2)$ matrix y_n at each site encoding the n^{th} instanton's orientation (for $N_f = 2$). We start with a completely random set of the initial y_n and make them evolve along the steepest descent of the net energy function until they converge to a local minimum. Then we repeat the process many times with different random sets of initial y_n . This gives us a short list of the local minima, from which we select the deepest — which is presumably the global minimum of the net energy function. This algorithm has convergence problems near the phase boundaries — where two or more local minima are almost degenerate — but it is good at finding all the orientation patterns that *might* be ground states for some combinations of background parameters and lattice geometries. And once we know all these patterns, we fit them all with common ansatz, calculate the net energy analytically as a function of the ansatz's parameters, and then we minimize it with a much higher precision to get the accurate phase diagram of the instanton lattice.

¹A Nahm transform can map the ADHM constraints to an electrostatic-like problem in 2D or 3D.

In this paper we put the instantons in a harmonic potential that rises at different rates along three out of four space coordinates; in terms of the 5D gauge coupling,

$$\frac{8\pi^2}{g_5^2(x_1, x_2, x_3, x_4)} = N_c \lambda M \times \left(1 + M_2^2 x_2^2 + M_3^2 x_3^2 + M_4^2 x_4^2 + O(M^4 x_{2,3,4}^4) \right) \quad (1.2)$$

with 3 independent parameters M_2, M_3, M_4 . Here we use different coordinate axes than in [1]: the 1D straight chain — or the long axis of the zigzag — runs along the x_1 axis instead of the x_4 . But the important difference is allowing for $M_3 \neq M_4$ and $M_2 \sim M_3 \sim M_4$ instead of $M_3 = M_4 = M' \gg M = M_2$; it is this difference which leads to the wide variety of instanton orientation patterns we shall see in this paper.

Specifically, we shall see that for the 1D instanton lattice in the almost-holographic background $M_2, M_3 \ll M_4$ (where M_2 and M_3 help to line up the instanton centers along the x_1 axis but do not affect the two-body forces between the instantons), there is a large degenerate family of ground states for the instanton orientations $y_n \in \text{SU}(2)/\mathbb{Z}_2$ (for $N_f = 2$). This family includes many periodic patterns in which all the y_n belong to a finite subgroup of the $\text{SU}(2)/\mathbb{Z}_2$: the anti-ferromagnetic pattern in which $y_{\text{odd } n}$ and $y_{\text{even } n}$ belong to 2 elements of a \mathbb{Z}_2 subgroup, the period = 4 pattern in which the y_n span the Klein group $\mathbb{Z}_2 \times \mathbb{Z}_2$, and two series of period = $2k$ patterns in which the orientations span a prismatic group $\mathbb{Z}_k \times \mathbb{Z}_2$ or a dihedral group D_{2k} . There are also many link-periodic patterns in which the relative orientations $y_n^\dagger y_{n+1}$ of neighboring instantons are periodic but the orientations y_n themselves are not periodic. But the vast majority of the degenerate ground states are not periodic at all.

Increasing the M_2 and M_3 parameters lifts the degeneracy. For $M_2 \ll M_3 \sim M_4$, the ground state of the 1D lattice is anti-ferromagnetic, while for $M_2 \sim M_3 \sim M_4$ the ground state is periodic with a longer period or link-periodic, depending on the M_2/M_3 ratio. Specifically, for $M_2 = M_3$ the instanton's orientation in the ground state span the Klein group, for other rational M_2^2/M_3^2 they span a dihedral group, while the irrational M_2^2/M_3^2 favor the link-periodic patterns.

For the zigzag-shaped chains of instantons we need $M_2 \ll M_3 \sim M_4$ to make sure the transition from a straight chain to a zigzag happens when lattice spacing \gg instanton radius — otherwise the two-body-force approximation would not work — but we may vary the M_3/M_4 . In this setting, the ground state of instanton orientations may follow one of four different patterns, depending on the zigzag amplitude and on the M_3/M_4 ratio. Besides the anti-ferromagnetic and the abelian $\phi < \pi$ patterns seen in [1], there are two non-abelian link-periodic pattern in which the $y_n^\dagger y_{n+1}$ “twists” for even and odd links of the lattice do not commute with each other. The non-abelian patterns appear only for $M_3 < M_4$, that's why we did not see them in [1].

Focusing on the “popcorn transition” from a straight chain of instantons to a zigzag-shaped chain (which is a two-layer lattice), we shall see that the thermodynamic order of this transition depends on the M_3/M_4 ratio: the transition is first-order for $M_3 < 0.725M_4$ but second order for $M_3 > 0.725M_4$. Likewise, when we further increase the instanton density — or rather, the 1D compression force which maintains it — the number and

the orders of the subsequent phase transitions between different orientation phases of the zigzag-shaped chain also depend on the M_3/M_4 ratio.

This paper is organized as follows: for the reader's convenience, in section 2 we review the basic features of holographic nuclear matter: effects of the large N_c and large λ limits, Sakai-Sugimoto-like models of holographic QCD, realization of holographic baryons as instantons of the flavor gauge symmetry, basic features of the holographic nuclear forces, and our basic approach to the multi-baryon systems and the bulk nuclear matter. In section 3, we argue that the interactions between multiple instantons include 2-body forces as well as irreducibly-3-body force, 4-body forces, etc., but when the distances between the baryons are much larger than the instantons' radii, the two-body forces dominate the net energy, while the 3-body, ..., forces are suppressed by powers of $(\text{radius}/\text{distance})^2 \ll 1$. First we give a heuristic argument, and then we back it up with a technical proof for $N_f = 2$; we also calculate the precise form of the net two-body force and its dependence on the two instantons' orientations.

In section 4 we use the two-body-force approximation to study 1D lattices of instantons. We start with the $M_2, M_3 \ll M_4$ background and show that there is a very large degenerate family of ground states of the instantons' orientations. Then we move on to the $M_2 \sim M_3 \sim M_4$ backgrounds; this changes the precise form of the two-body force in a way that lifts the degeneracy between the orientation states. Consequently, the ground state of the lattice has a specific periodic or link-periodic pattern of orientations, depending on the M_2/M_3 ratio.

In section 5 we study zigzag-shaped instanton chains in the $M_2 \ll M_3 \sim M_4$ backgrounds. After a few preliminaries, we use numerical methods to find the lowest-energy state of instanton orientations without presuming any particular symmetries. We repeat the calculation for 50 different lattice geometries ($\epsilon/D = 0.01, \dots, 0.99$ by 0.02) and 50 values of the M_3/M_4 ratio; the results are presented on the phase diagram 4 on page 48. Our numeric code is not very accurate and has convergence problems near the phase boundaries, but it is good at identifying the symmetries of the lowest-energy patterns. Using these symmetries, we construct a 3-parameter ansatz that includes all the preferred patterns, then calculate the net energy as analytic function of the parameters and minimize it. For the reader's convenience, the energy calculation and minimization is described in the appendix, while the resulting phase diagram is presented in figure 5 (page 52) in the main body of section 5.

In the last part of section 5 we treat the zigzag amplitude and the lattice spacing as dynamic parameters, vary the compression force of the 1D instanton lattice, and study the popcorn transitions from the straight chain to the zigzag and between different orientations phases of the zigzag. We show that the net number of such transitions — 1, 2, or 3 — and their orders depend on the M_3/M_4 ratio. The overall phase diagram in the compression versus M_3/M_4 plane is presented in figure 6(a) on page 56; we also present the phase diagrams in the chemical potential versus M_3/M_4 plane (figure 6(b) on page 56) and in the instanton density versus M_3/M_4 plane (figure 7 on page 57).

Finally, in section 6 we summarize our results and give a short list of open questions we hope to address in the near future.

2 Review of holographic nuclear physics

In this section we review the basics of holographic nuclear physics that we shall need later in this paper. The *holography* here means gauge-gravity duality, and the semiclassical description of the gravity side of the duality requires the limits of large N_c and large 't Hooft coupling $\lambda = g_{\text{YM}}^2 \times N_c$ [20]. In the first two subsections 2.1, 2.2 we discuss the general consequences of these limits for the nuclear physics. In subsection 2.3 we get more specific and review the Sakai-Sugimoto-like models of holographic QCD. There are many such models which differ in countless details — and none of them is exactly dual to the QCD, — but their basic features are rather similar, and the Sakai-Sugimoto model serves as a prototype. In subsection 2.4 we introduce the baryons and explain their holographic duality to the instantons of the flavor gauge symmetry. We also review the interactions and the self-interactions of such instantons in the large λ limit and explain why they have small radii $a \propto \lambda^{-1/2}$. Finally, in subsection 2.5 we address the multi-instanton systems dual to nuclear matter and explain how the interactions between the baryons appear in at the first order in the $1/\lambda$ expansion. We also explain the technical difficulties in handling infinite crystals of baryons and preview a shortcut around those difficulties that we shall take in section 3.

2.1 Nuclear matter in the large N_c limit

QCD perturbation theory in the large N_c limit is dominated by planar gluon diagrams while contributions of the non-planar diagrams and of the quark loops are suppressed. At zero temperature and chemical potential below the mass of the baryon, the large N_c QCD is a theory of free mesons and glueballs whose interactions are suppressed by powers of N_c (except for interactions with the baryons). The masses of mesons and glueballs are $O(\Lambda_{\text{QCD}})$, and they scale as N_c^0 . On the other hand, the baryons — which are made out of N_c quarks — have masses of order $N_c \times \Lambda_{\text{QCD}}$, so their relative abundance in thermal equilibrium is exponentially suppressed. The interaction energy between baryons also scales as N_c .

The $T - \mu$ phase diagram of the large- N_c QCD is believed to look like figure 1 on the next page. (For a review see [19] and references therein, also [29].) We assume the 't Hooft's $N_c \rightarrow \infty$ limit in which the number of flavors remains finite. Consequently, the dynamics of the theory is dominated by the gluons, the quarks are sensitive to the gluonic background, but the backreaction from the quarks to the gluons is suppressed by $N_f/N_c \ll 1$. Thus, at lower temperatures there is confinement but for increasing temperature there is a first order transition to the deconfined phase. The transition temperature T_d is at the Λ_{QCD} scale and it's almost independent of the quark chemical potential $\mu_q = \mu_b/N_c$ (as long as μ_q does not grow with the N_c).

It is not clear whether for $N_c \rightarrow \infty$ the deconfining phase transition coincides with the chiral symmetry restoration for the light quarks. Several field theory arguments — for example [19, 21, 22] — suggest that for $\mu_q = 0$ the two transitions should happen at the same point. However, these arguments do not work for $\mu_q > 0$ [19], and there are other arguments for the existence of confined but chirally restored phases and/or deconfined

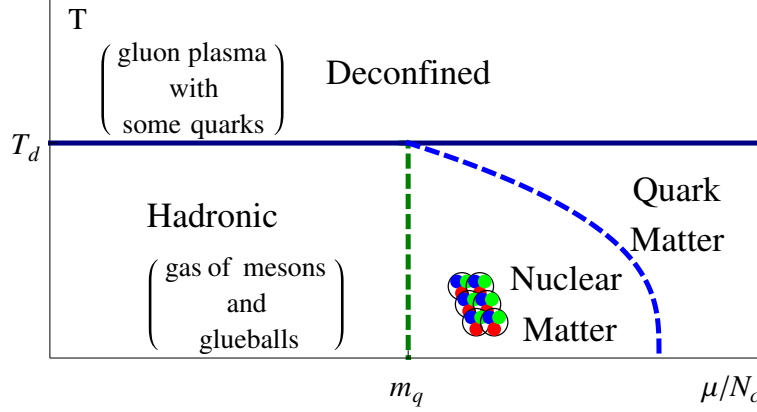


Figure 1. Conjectured phase diagram for QCD at large N_c [19].

phases where the chiral symmetry remains broken. Moreover, some holographic models — for example [4] — have deconfined but chirally broken phases even at zero chemical potentials.

For temperatures below the deconfining transition t_d and baryon chemical potentials below the baryon mass — or equivalently for $\mu_q \lesssim m_q \equiv M_b/N_c$ — the thermal state of the theory is a gas of glueballs and mesons with almost no baryons or antibaryons. But at $\mu_q \approx m_q$ there is an abrupt phase transition to the bulk nuclear matter with finite baryon density. But unlike the ordinary nuclear matter which is in the quantum liquid state, the large N_c nuclear matter is crystalline solid since the ratio of kinetic energy to potential energy decreases with N_c . Indeed, the potential energy of baryon-baryon forces scales like N_c ; more precisely [23], in the large N_c limit the two-baryon potential becomes

$$V \sim N_c \times A_C(r) + N_c \times A_S(r) (\mathbf{I}_1 \cdot \mathbf{I}_2) (\mathbf{J}_1 \cdot \mathbf{J}_2) + N_c \times A_T(r) (\mathbf{I}_1 \cdot \mathbf{I}_2) (3(\mathbf{n} \cdot \mathbf{J}_1)(\mathbf{n} \cdot \mathbf{J}_2) - (\mathbf{J}_1 \cdot \mathbf{J}_2)) + O(1/N_c), \quad (2.1)$$

for some N_c -independent profiles $A_C(r)$, $A_S(r)$, and $A_T(r)$ for the central, spin-spin, and tensor forces; their overall magnitudes are $A \sim \Lambda_{\text{QCD}}$ for $r \sim 1/\Lambda_{\text{QCD}}$. Classically, this potential tries to organize the baryons into some kind of a crystal where the distances between neighboring baryons do not depend on the N_c while the binding energy (per baryon) scales like $N_c \times O(\Lambda_{\text{QCD}})$. In quantum mechanics, the baryons in such a crystal behave like atoms in ordinary crystals — they oscillate in their potential wells with zero-point kinetic energies

$$K \sim \frac{\pi}{2m_B d^2} \sim \frac{\Lambda_{\text{QCD}}}{N_c} \frac{1}{d^2}, \quad (2.2)$$

where $d \sim 1/\Lambda_{\text{QCD}}$ is the N_c -independent diameter of the potential well. Therefore, at zero temperature the ratio of kinetic energy to the potential energy scales like

$$\frac{K}{V} \sim \frac{1}{N_c^2} \quad (2.3)$$

and becomes very small for large N_c . At higher temperatures the kinetic energies of baryons become larger, $K \sim T$, but in the confined phase we are limited to $T < T_d \sim \Lambda_{\text{QCD}}$ and

hence $K < O(\Lambda_{\text{QCD}})$. Consequently, the kinetic to potential energy ratio scales like

$$\frac{K}{V} \sim \frac{1}{N_c} \quad (2.4)$$

— which is larger than (2.3) but still becomes small in the large N_c limit. Consequently, for large N_c neither zero-point quantum motion nor thermal motion of baryons can melt the baryon crystal, so nuclear matter remains solid all the way to the deconfining temperature.

Before holography, the best models for the large- N_c nuclear crystals were lattices of skyrmions. In this framework, Igor Klebanov [24] had found a curious phase transition from a lattice of whole skyrmions at low chemical potentials to a denser lattice of half-skyrmions at higher potentials. (According to [25, 26], the whole-skyrmion lattice at low potentials has FCC geometry while the half-skyrmion lattice at higher potentials is simple cubic.) In the half-skyrmion-lattice phase, the order parameter for the chiral symmetry breaking vanishes after space averaging, so in QCD terms the transition is interpreted as chiral symmetry restoration at high μ_q .

In QCD with $N_c = 3$ and two massless flavors there is a similar chirally-symmetric phase at low T and high μ_q . This phase is a quark liquid rather than a baryon liquid (the quarks are no longer confined to individual baryons) and there is a condensate of quark pairs making this liquid a color superconductor. But for large N_c the situation is more complicated: there is no color superconductivity, and there is no deconfinement for $T < T_d$. Instead, the dense cold nuclear matter forms a phase which combines the features of the baryonic and quark phases: the quarks fill up a Fermi sea, but the interactions near the Fermi surface are strong, so the excitations are not free quarks or holes but rather meson-like quark-hole pairs or baryon-like states of N_c quarks. MacLerran et al. have dubbed this phase “quarkyonic”.

For $\mu_q \gg \Lambda_{\text{QCD}}$, the interior of the Fermi sea is chirally symmetric, but near the Fermi surface the symmetry is broken by the chiral density waves [27]. (Although Son and Shuster [38] argue that such waves develop only for very large $N_c > 10^3$.) To be precise, the chiral density waves mix the chiral symmetry of the quarkyonic phase with the translational symmetry rather than simply break it. Averaging over space restores the chiral symmetry, just like it happens for the lattice of half-skyrmions.

On the other hand, for μ_q just above $m_q = M_b/N_c$, the baryonic crystal has a completely broken chiral symmetry. Thus, at some critical $\mu_q^{(c)} = m_q \times \text{a few}$, there should be a chiral symmetry restoring phase transition from the baryonic crystal to a distinct quarkyonic phase. In holography, this transition is believed to be dual to the “popcorn transition” from a 3D instanton lattice to a 4D lattice.

2.2 Effect of the large λ limit on the holographic nuclear matter

In holography, the semiclassical description of the gravity side of the gauge-gravity duality in terms of metric, fluxes, and branes requires the limits of large N_c and also large 't Hooft coupling $\lambda = N_c g_{\text{YM}}^2$. In the large λ limit, the baryons become very heavy: in units of the mesonic mass scale $M \sim \Lambda_{\text{QCD}}$, the baryon mass is $M_b \sim \lambda N_c M$. However, the interactions between the baryons do not grow with λ : even for two baryons right on top of each other,

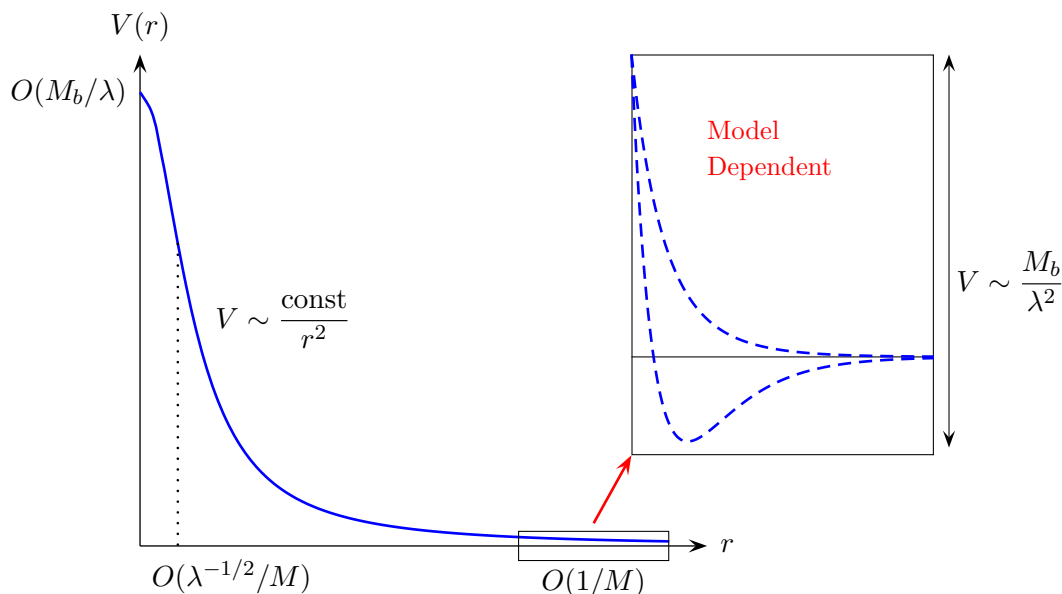


Figure 2. Two-body nuclear potential in holographic QCD.

the repulsive potential between them is only $V \sim N_c M \sim M_b/\lambda$. At larger distances, the forces are even weaker since the hard-core radius of a holographic baryon shrinks with λ as $R_b \sim M^{-1}\lambda^{-1/2}$. Outside this radius, the repulsive potential decreases as $1/r^2$ until $r \sim M^{-1}$, at which point it becomes dependent on the meson mass spectrum of a specific holographic model: in some models, the potential becomes attractive for $r \gtrsim M^{-1}$ while in others it remains repulsive at all distances. The overall picture is shown in figure 2.

Since the nuclear forces are so weak in the holographic QCD, all transitions between different phases of the cold nuclear matter happen at chemical potentials μ_b very close to the baryon mass: just below M_b we have glueball/meson gas (or vacuum for $T = 0$), while just above M_b we have dense quark matter. To see the baryonic matter phase (or any other intermediate phases) we need to zoom into the $\mu_b \approx M_b$ region. Figure 3 (on the next page) illustrates this point: to see the baryonic-matter phase between the vacuum and the quark-matter phases on the plot of baryon density ρ as a function of the chemical potential μ , we need to zoom into the narrow range $\mu - M_b = O(M_b/\lambda)$. The figure also shows that the thermodynamic order of the phase transition between the vacuum (or the meson/glueball gas for $T > 0$) and the baryonic matter depends on the sign of the long-distances nuclear force. If the force becomes attractive at long distances, then bulk baryonic matter exists at zero external pressure and has $\mu = M_b$ – binding energy. Consequently, the transition from the vacuum (or gas) to the nuclear matter (in the form of a baryonic crystal) is first-order as shown on figure 3(a). On the other hand, if the nuclear forces are repulsive at all distances, then the bulk nuclear matter does not exist except at positive external pressures, and its chemical potential must be $\mu > M_b$. Moreover, μ raises monotonically with the pressure and the density, so the transition from the vacuum to the bulk nuclear matter is second-order as shown on figure 3(b).

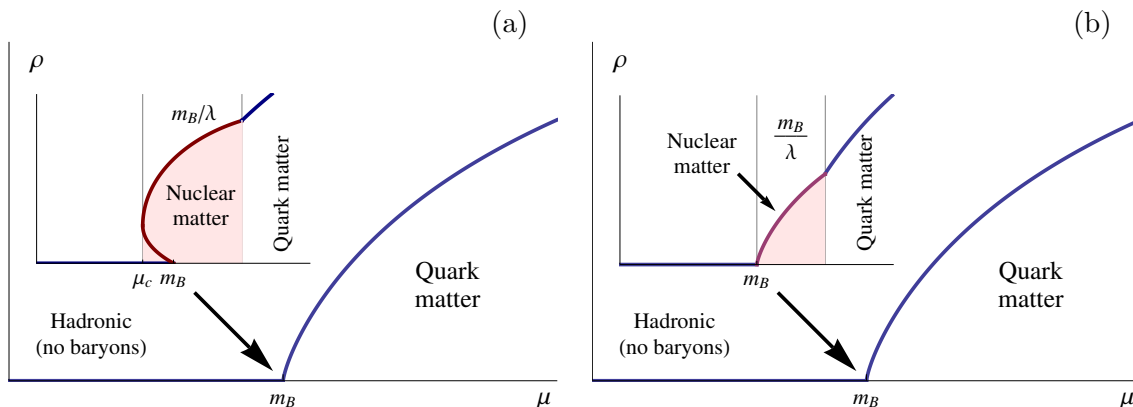


Figure 3. Density as a function of the chemical potential in the large λ regime for attracting (a) and repelling (b) baryons. The nuclear matter phase is confined to a narrow window of the order $\Delta\mu \sim B_B/\lambda$. In the naive diagram the transition occurs directly from the no-baryon to a quark phase.

In this article we do not wish to focus on a particular model of holographic QCD, so we make no assumptions about the long-distance nuclear forces. Consequently, we cannot say anything specific about the transition from the vacuum to the baryonic phase. Instead, we focus on the transition from the baryonic phase to the quarkyonic phase, which correspond in the holographic picture to changing the lattice geometry of the instanton crystal, from a 3D lattice, through a sequence of intermediate steps, to a 4D lattice. Or rather, that’s the goal of our program; in this article we shall focus on a simplified problem, namely the transitions between 1D and 2D instanton lattices.

2.3 Sakai-Sugimoto-like models of holographic QCD

Some gauge theories — especially the $\mathcal{N} = 4$ SYM — have exact holographic duals, where both sides of the duality follow as IR limits of the same string-theoretical construction, and all the extra degrees of freedom are superheavy. Alas, the Quantum Chromo-Dynamics — or even the large N_c limit of QCD — is not so lucky: it either does not have an exact holographic limit, or it has not been found yet. Instead, there is a large number of “holographic QCD” *models*, which are dual not to the QCD as such but rather to some QCD-like theories with a lot of extra junk that the real QCD does not have. Such models make for “spherical horse in a vacuum” approximations — good enough to get qualitative understanding, but not too accurate to make useful numerical predictions.

In this article we are interested in a particular class of holographic QCD models exemplified by the Sakai-Sugimoto model [3]. Here are some general features of such models:

- The construction starts with N_c coincident D-branes, which span the Minkowski space \times a compact cycle. Typically, one uses D3 branes at a singularity, or D4-branes wrapping a circle, although some constructions use D5 branes on a 2-cycle, etc. For weak ’t Hooft couplings $\lambda = N_c \times \alpha_{\text{QCD}} \sim N_c \times g_{\text{str}} \ll 1$, the open strings between the branes give rise to the gluons of the $U(N_c)$ gauge theory (plus a lot of extra junk).

- For $\lambda \gg 1$, the D-branes merge into a big fat black brane. All we see outside the horizon is a warped space-time geometry and the Ramond-Ramond fluxes induced by the conserved charges of the D-branes. Modes of the bulk-field excitations in this geometry are dual to the QCD *glueballs*. Or rather, some modes are dual to the glueballs while other models are dual to the extra junk that the real QCD does not have 😞
- To make the hadrons (mesons and baryons), one adds N_f flavor D-branes, usually D5, D6, D7, or D8. At weak λ , the open strings connecting the color branes to the flavor branes give rise to the quarks and the anti-quarks.
- In the holographic limit $N_c \rightarrow \infty$ and $\lambda \rightarrow \infty$, but the N_f stays finite and the $N_f \times g_{\text{str}}$ remains weak. Consequently, the flavor branes remain D-branes (rather than merge into a black brane of their own), but now they live in a warped geometry with fluxes induced by the N_c color branes.
- The open strings between the flavor branes give rise to N_f^2 vector and scalar fields living on those branes. The 4D modes of these vector and scalar fields are dual to the QCD *mesons*.
- The YM instantons of the vector fields on the flavor branes are dual to the QCD *baryons*, see subsection 2.4 for more details.

The type specimen of this class — the Sakai-Sugimoto model [3] — was reviewed in detail in our previous paper [1] (as well as by countless other authors), so here we may be brief. In the $\lambda \gg 1$ limit, the bulk geometry of the model is a warped product $\mathbf{R}_{\text{Minkowski}}^{1,3} \times S^4 \times \text{cigar}$, while the flavor degrees of freedom live on D8 branes spanning $\mathbf{R}^{1,3} \times S^4 \times$ a U-shaped line on the cigar. The effective action for the flavor fields is dominated by the Dirac-Born-Infeld and Chern-Simons terms,

$$S = S_{\text{DBI}} + S_{\text{CS}} = T_8 \int_{D8} d^9x \text{Str} \left(\sqrt{-\det(g_{MN}[\phi] + 2\pi\alpha' \mathcal{F}_{MN})} \right) + T_8 \int_{D8} C_3 \wedge \exp(2\pi\alpha' \mathcal{F}). \quad (2.5)$$

where T_8 is the D*-brane tension, $g_{MN}[\phi]$ is the induced 9D metric on the brane perturbed by the scalar fields ϕ , \mathcal{F}_{MN} are the $U(N_f)$ gauge fields, and Str denotes the symmetrized trace.

For our purposes, we are interested in the flavor field configurations which do not depend on the S^4 coordinates, while the vector fields do not have any components along the S^4 . Therefore, we dimensionally reduce the flavor theory to just five dimensions — the Minkowski coordinates $x^{0,1,2,3}$, and the z coordinate of the U-shaped line on the cigar. The effective 5D theory has vector and scalar fields in the adjoint multiplets of $U(N_f)$ with an effective action

$$S_{\text{eff}} = \int d^4x dz \left[\begin{aligned} & \frac{1}{g_{5d}^2(z)} \text{tr} \left(-\frac{1}{2} \mathcal{F}_{mn} \mathcal{F}^{mn} + D_m \Phi D^m \Phi \right) - V(\Phi, z) \\ & + N_c \times \omega_{\text{CS}}(\mathcal{A}, \mathcal{F}) + \frac{N_c C(z)}{16\pi^2} \times \text{tr}(\Phi \mathcal{F}_{mn} \mathcal{F}^{mn}) \\ & + \text{higher-derivative terms} \end{aligned} \right]. \quad (2.6)$$

In our notations the z coordinate is normalized such that the background 5D metric is conformal, $ds^2 = A(z) \times (dt^2 - d\vec{x}^2 - dz^2)$, the scalar fields Φ are normalized so they have a similar kinetic energy coefficient $1/g_{5d}^2(z)$ to the vector fields, while the scalar potential has general form

$$V(\Phi, z) = m_\phi^2(z) \times \text{tr}(\Phi^2) + p(z) \times \text{tr}(\Phi^3) + \dots$$

but its details are not important for our purposes.

For other holographic QCD models in our class, the flavor branes have different dimensionalities and span different compact cycles, but only five of the dimensions are interesting, the rest giving rise to junk mesons without QCD analogues. When we dimensionally reduce such models to five dimensions, they end up with effective actions similar to (2.6), the only difference being in the z -dependent $g_{5d}^2(z)$ and $C(z)$ couplings. In general, these couplings depend on z on the Kaluza-Klein scale of the theory, while in this article we are going to be interested at the much shorter distance scale of the baryon radius $R_b \sim 1/(\lambda^{1/2} M_{KK}) \ll 1/M_{KK}$, so all we need is the expansion

$$\frac{8\pi^2}{g_{5d}^2(z)} = N_c \lambda M_{KK} (B + D M_{KK}^2 \times z^2 + O(M_{KK}^4 z^4)) \quad (2.7)$$

for some numeric constants B and D ; for the antipodal Sakai-Sugimoto model $B = 1/27\pi$ and $D = 1/3$, while other models have different values. Likewise, we may approximate $C(Z) \approx C = \text{const.}$

The baryons are holographically dual to instantons of the flavor gauge fields, so the multi-baryon system have strong self-dual $SU(N_f)$ magnetic fields, $\mathcal{F}_{\mu\nu} = \frac{1}{2} \epsilon_{\mu\nu\rho\sigma} \mathcal{F}^{\rho\sigma}$ ($\mu, \nu = 1, 2, 3, z$). For such self-dual fields, the DBI action reduces to Yang-Mills, so we may neglect the higher-derivative terms in the action (2.6). At the same time, the Chern-Simons term

$$N_c \times \omega_{\text{CS}}(\mathcal{A}, \mathcal{F}) = \frac{N_c}{24\pi^2} \times \text{tr} \left(\mathcal{A} \wedge \mathcal{F}^2 - \frac{i}{2} \mathcal{A}^3 \wedge \mathcal{F} - \frac{1}{10} \mathcal{A}^5 \right) \quad (2.8)$$

turns the instanton number density

$$I(x, z) = \frac{\epsilon^{\kappa\lambda\mu\nu} F_{\kappa\lambda}^a F_{\mu\nu}^a}{32\pi^2} \quad (2.9)$$

into a charge density which excites the Coulomb electric fields. For $N_f = 2$ such electric fields are purely abelian; indeed, splitting the gauge fields \mathcal{F}_{mn} into $SU(2)$ fields F_{mn} and $U(1)$ fields \hat{F}_{mn} , we have

$$\mathcal{L}_{\text{CS}} = \frac{N_c}{16\pi^2} \hat{A} \wedge \text{tr}(F^2) + \frac{N_c}{96\pi^2} \hat{A} \wedge \hat{F}^2 = \hat{A}^0(x, z) \times \frac{N_c}{2} I(x, z) + \dots, \quad (2.10)$$

so an $SU(2)$ instanton has effective $U(1)$ electric charge $+\frac{N_c}{2}$. For $N_f > 2$, the instantons have both abelian and nonabelian electric charges, which we shall explain in the next section 2.4. In any case, the electric fields lead to Coulomb repulsive forces between the holographic baryons; for distances larger than the baryon radius but smaller than the Kaluza-Klein scale, the repulsive potential behaves as $1/r^2$.

Besides repulsive forces, most holographic models also have attractive forces due to backreaction of the strong magnetic fields on the geometry of the flavor brane. In the effective 5D action (2.6), this backreaction is encoded in the scalar-vector coupling $\text{tr}(\Phi\mathcal{F}\mathcal{F})$ — the strong magnetic fields act as sources of scalar fields, which lead to the attractive forces. For the self-dual magnetic fields, the scalar fields couple to exactly the same source as the electric fields A^0 , except for the overall coefficient $C(z) \approx C$, so at short and intermediate distances between the baryons the attractive and repulsive forces have constant ratio

$$\frac{\text{attractive}}{\text{repulsive}} = C^2(z) \approx C^2. \quad (2.11)$$

In the antipodal Sakai-Sugimoto model, the leading-order backreaction vanishes due to a geometric symmetry, hence $C = 0$ and there are no attractive forces at all. But other models — and even the non-antipodal versions of the Sakai-Sugimoto model [2] — have $C > 0$ and hence non-trivial attractive forces. The direction of net the short-distance force depends on the sign of $1 - C^2$; in this paper, we assume $1 - C^2 > 0$, so at short distances there is a repulsive hard core.

2.4 Baryons as flavor instantons

In the old hadronic string, the baryons were made out of Y-shaped configurations of 3 open strings connected to each other at one end; the other end of each string is connected to a quark. To realize this picture in holographic QCD, we need a *baryon vertex* (BV) — some object connected to N_c open strings. The other ends of the strings are connected to the flavor branes and act like the quarks; this gives the baryon its flavor quantum numbers. Witten had constructed the baryon vertex for the $\text{AdS}_5 \times S^5$ model from a D5 brane wrapping the S^5 sphere [30]; the generalized versions of this construction [31–33] use D_p branes wrapping compact cycles carrying $O(N_c)$ Ramond-Ramond fluxes.

In the Sakai-Sugimoto version of this construction, the baryon vertex is realized as a D4 brane wrapped around the S^4 sphere (but localized in all other dimensions except the time). The S^4 carries N_c units of the F_4 Ramond-Ramond flux,

$$\frac{1}{(2\pi)^3 l_s^3} \int_{S^4} F_4 = N_c, \quad (2.12)$$

so the Chern-Simons coupling of this flux to the $U(1)$ gauge field \mathcal{B} living on the D4 brane acts a N_c units of the net electric charge for the B_0 :

$$T_4 \int_{D4} C_3 \wedge e^{2\pi\alpha' dB} = N_c \int_{\text{time}} B_0 dt. \quad (2.13)$$

In a compact space like S^4 , the net electric charge must vanish. To cancel the charge (2.13) we need to connect the D4 brane to open strings. The back end of an oriented open string has electric charge -1 , so we must connect the D4 brane with N_c such strings; their front ends connect to the D8 flavor branes (since the strings do not have any other place to end) and act as N_c quarks.

We may put the D4 brane anywhere in space and anywhere on the cigar. However, the S^4 volume increases with the u coordinate, so the lowest-energy location of the D4 is

the cigar's tip $u = u_\Lambda$. At other location, the brane feels a gravity-like force pulling it down to the tip. However, the strings connected to the baryon vertex pull it towards the flavor D8 branes; in the non-antipodal models the D8 branes do not reach the cigar's tip, so the strings pull the baryonic vertex up from the tip towards the lowest point u_0 of the flavor branes. The ‘tag of war’ between the upward and downward forces on the baryon vertex determines its ultimate location. In some models, the forces reach equilibrium for the baryon vertex hanging on strings below the flavor branes [52], while in many other models — including the non-antipodal Sakai-Sugimoto — the string forces win and pull the baryon vertex all the way to the lowest point u_0 of the flavor branes [40].

In all such models, the baryonic vertex is a D_p brane completely embedded in a stack of D_{p+4} flavor branes, so it is equivalent to zero-radius Yang-Mills instanton of the $U(N_f)$ gauge symmetry on the flavor branes, and for $N_f > 1$ it may be smoothly inflated to a finite-radius instanton [41–45]. In $p + 5$ dimensions of the flavor branes, this instanton is a fat D_p -brane wrapping some compact cycle, but once we dimensionally reduce to 5 dimensions, the instanton becomes a finite-size particle. Thus, *in the low-energy effective 5D theory of the holographic QCD, a baryon is realized as a finite-size instanton of the $U(N_f)$ gauge theory.*

Our research program of studying the instanton crystals as holographic duals of the cold nuclear matter stems from this realization of baryons as instantons in 5D, but it does not depend on any specific details of the Sakai-Sugimoto model. We use that model as a prototypical example, but any other model where the baryons are realized as instantons would be just as good for our purposes, for example the model of [34] for some 7-brane geometries [52], or the $AdS_5 \times S^1$ model of [46] (the baryons of that model are studied in section 6 of [40]).

In the baryon-vertex picture, each of the N_c strings connecting the vertex to the flavor branes has electric charge $1/N_f$ under the abelian $U(1)$ subgroup of the $U(N_f)$, so the whole baryon has abelian charge N_c/N_f . In the instanton picture, the same electric charge obtains from the Chern-Simons coupling between the abelian electric field and the non-abelian magnetic fields of the instanton.

$$\begin{aligned} S_{CS} &= \frac{N_c}{24\pi^2} \int tr \left(\mathcal{A} \mathcal{F}^2 - \frac{i}{2} \mathcal{A}^3 \mathcal{F} - \frac{1}{10} \mathcal{A}^5 \right) \supset \frac{N_c}{N_f} \int \hat{A}_{U(1)} \wedge \frac{\text{tr}(F \wedge F)_{SU(N_f)}}{8\pi^2}, \\ &\supset \frac{N_c}{N_f} \int d^5x \hat{A}_0(x) \times I(x) \end{aligned} \quad (2.14)$$

where $I(x)$ is the instanton number density of the magnetic fields. For $N_f \geq 3$, the Chern-Simons couplings also endow instantons with non-abelian electric charges. Altogether,

$$S_{CS} \supset \int d^5x A_0^a(x) \times Q_{el}^a I(x) \quad (2.15)$$

where the net electric charge Q_{el}^a has both abelian and non-abelian components; in $N_f \times N_f$

matrix language,

$$\text{for } F_{\text{magnetic}}^{\mu\nu} \in \left(\begin{array}{c|c} \text{SU}(2) & \mathbf{0} \\ \hline \mathbf{0} & \mathbf{0} \end{array} \right), \quad Q_{\text{el}} = \frac{N_c}{2} \times \left(\begin{array}{c|c} \mathbf{1} & \mathbf{0} \\ \hline \mathbf{0} & \mathbf{0} \end{array} \right), \quad \text{tr}(Q_{\text{el}}^2) = \frac{N_c^2}{2}. \quad (2.16)$$

Abelian or nonabelian, the like-sign electric charges repel each other; it is this Coulomb repulsion between different parts of the same instantons that prevents it from collapsing to a point-like D-brane.

However, the instantons do not grow large because the 5D gauge coupling decreases away from the $z = 0$ hyperplane: a large instanton would spread into regions of space where the coupling is weaker, and that would increase the instanton's energy. Instead, the equilibrium radius of the instanton scales like

$$a_{\text{instanton}} \sim \frac{1}{M_{\text{KK}} \times \sqrt{\lambda}}. \quad (2.17)$$

For a holographic model of a baryon, this radius is unrealistically small. Indeed, using the ρ meson's mass as a unit, the real-life baryon radius $R_b \sim 3.4 M_\rho^{-1}$, while in holography $a \ll M_\rho^{-1} \sim M_{\text{KK}}^{-1}$. Moreover, it raises the question of whether we may adequately describe so small an instanton using the DBI + CS action, or perhaps we need to include the higher-order stringy corrections.² On the other hand, assuming the DBI + CS description is OK, the radius as in eq. (2.17) allows for consistent expansion of the instanton's parameters in powers of λ^{-1} [5]. In particular, the leading contribution to the instanton's mass is $O(\lambda \times N_c M_{\text{KK}})$ while the corrections due to z -dependent gauge coupling and due to Coulomb self-repulsion are both $O(N_c M_{\text{KK}}) \sim M_I/\lambda$.

To see how that works, consider a static instanton — a time-independent configuration of $\text{SU}(N_f)$ magnetic fields, plus the electric fields induced by the CS couplings, and the scalar fields induced by the $\text{tr}(\Phi FF)$ coupling to the magnetic fields. Since the canonically normalized couplings in 5D are $O(\lambda^{-1/2})$, the leading contribution to the instanton's energy comes from the magnetic field. In the DBI approximation³

$$E_{\text{DBI}} = \int d^3\mathbf{x} \int dz \frac{1}{2g_{\text{YM}}^2(z)} \times \text{Str} \left(\sqrt{\det(K(z)\delta_{\mu\nu} + F_{\mu\nu})} - K^2(z) \right) \quad (2.18)$$

²In string theory, the DBI + CS action for the gauge fields on a D-brane is exact for *constant* tension fields \mathcal{F}_{mn} , however strong. But for the variable tension fields, the DBI action includes all powers of the \mathcal{F}_{mn} but neglects their derivatives $D_k \mathcal{F}_{mn}$, $D_k D_p \mathcal{F}_{mn}$, \dots . It is not clear what effect (if any) such higher-derivative terms would have on a small instanton. In a supersymmetric background, the instanton is BPS and its net mass is protected against stringy corrections, so the DBI action — or even the Yang-Mills action — gives the exact value. But what happens to small instanton in non-supersymmetric backgrounds is an open question.

³In string theory, the complete action for scalar and vector fields living on a D-brane starts with the DBI and CS terms but also includes an infinite number of terms involving $D_n \mathcal{F}_{mn}$ and the higher derivatives. The DBI + CS action is merely a long-distance approximation that allows for strong gauge field tensions \mathcal{F}_{mn} as long as their gradients are small. It is not clear how well this approximation works for small instantons, but at present we do not have a better approximation.

where $K(z) = (2\pi\alpha')^{-1} \times g_{11}(z) \sim 1/(\lambda M_{KK}^2 \ell_s^4)$. Both $g_{YM}(z)$ and $K(z)$ depend on the z coordinate on the distance scale $O(1/M_{KK})$, so for instantons of much smaller size we may start with the approximation of constant K and constant 5D gauge coupling. And in this approximation, the DBI energy is minimized by the magnetic fields that are exactly self-dual (in the 4D space of (x^1, x^2, x^3, z)); moreover, the DBI energy of an instanton is equal to its Yang-Mills energy

$$E_{\text{DBI}}(\text{instanton}) = E_{\text{YM}}(\text{instanton}) = \frac{8\pi^2}{g_{\text{YM}}^2} \quad (2.19)$$

regardless of its radius and of the K parameter of the DBI action.

To determine the equilibrium radius of an instanton we need to calculate its energy to the next order in $1/\lambda$ expansion. To this order, we assume the magnetic fields to be exactly self-dual — which allows us to use the YM action instead of DBI even for small instantons — but the gauge coupling is z -dependent, and we also account for the electric and the scalar fields. For small instantons, we may approximate the 5D gauge coupling as

$$\frac{8\pi^2}{g^2(z)} = N_c \lambda M_{KK} (B + D M_{KK}^2 \times z^2 + O(M_{KK}^4 z^4)) \quad (2.20)$$

for some numerical constants B and D ; for the Sakai-Sugimoto model $B = \zeta/27\pi$ and $D = (8\zeta^3 - 5)/9\zeta^2$, while other models may have different values. Consequently, the YM energy of the nonabelian magnetic fields evaluates to

$$E_{\text{NA}} = N_c \lambda M_{KK} \times \left(B + D M_{KK}^2 \times Z^2 + D M_{KK}^2 \times \frac{a^2}{2} \right) \quad (2.21)$$

where a is the instanton's radius and Z is the z coordinate of its center.

The electric potentials A_0^a couple to the $F_{\mu\nu} \tilde{F}^{\mu\nu}$ products of the magnetic fields while the scalar potentials Φ^a couple to the $F_{\mu\nu} F^{\mu\nu}$. For the self-dual magnetic fields, both potentials couple to the same source $I(\mathbf{x}, z) \times Q_{\text{el}}^a$, and the only difference is the coupling strength ratio $C(z)$ (cf. eqs. (2.6) and (2.11)). For small instantons we may neglect the z -dependence of this ratio and let $C(z) \approx C(0) \equiv C$, and as long as we do not go very far from the instanton (for distances $r \ll R_{KK}$), we may also neglect the z -dependence of the gauge coupling. Consequently, both electric and scalar potentials become the 4D Coulomb potentials

$$\Phi(\mathbf{x}, z) = C \times A_0(\mathbf{x}, z) = \frac{2Q_{\text{electric}}}{BN_c \lambda M_{KK}} \int d^3 \mathbf{x}' \int dz' \frac{I(\mathbf{x}', z')}{(\mathbf{x} - \mathbf{x}')^2 + (z - z')^2}. \quad (2.22)$$

The electric fields lead to repulsive forces between the charges while the scalar forces lead to the attractive forces. Altogether, the net Coulomb energy amounts to

$$\begin{aligned} E_C &= \frac{B}{8\pi^2} N_c \lambda M_{KK} \times \int d^3 \mathbf{x} dz \text{tr}((\nabla A_0)^2 - (\nabla \Phi)^2) \\ &= (1 - C^2) \times \frac{N_c}{4B\lambda M} \int d^3 \mathbf{x} d^3 \mathbf{x}' dz dz' \frac{I(\mathbf{x}, z) \times I(\mathbf{x}', z')}{(\mathbf{x} - \mathbf{x}')^2 + (z - z')^2}, \end{aligned} \quad (2.23)$$

and for a single instanton of radius a it evaluates to

$$E_C = \frac{1 - C^2}{B} \times \frac{N_c}{5\lambda M_{\text{KK}} a^2}. \quad (2.24)$$

Note: for $C < 1$ the electric fields are stronger than the scalar fields and the net Coulomb energy of the instanton is positive — which makes for a net self-repulsive force that prevents the instanton from shrinking to zero radius. In models with $C > 1$ (assuming they exist), the scalar fields would be stronger, the net Coulomb energy would be negative, which means a self-attractive force rather than self-repulsive. In such a model, the instanton would shrink to zero radius and our approximations would not be valid.

Altogether, the net energy of an instanton amounts to

$$E(\text{instanton}) = N_c \lambda M_{\text{KK}} \times \left(B + D M_{\text{KK}}^2 \times Z^2 + D M_{\text{KK}}^2 \times \frac{a^2}{2} + \frac{1 - C^2}{5B} \times \frac{1}{\lambda^2 M_{\text{KK}}^2 a^2} \right) \quad (2.25)$$

Note that for $a^2 \sim 1/(\lambda M_{\text{KK}}^2)$ both radius-dependent terms here are $O(1/\lambda)$ corrections to the leading term. Minimizing the net energy, we find the equilibrium value of the instanton radius

$$a = \frac{1}{M_{\text{KK}} \sqrt{\lambda}} \times \sqrt[4]{\frac{2(1 - C^2)}{5BD}}, \quad (2.26)$$

the instanton center is in equilibrium at $Z = 0$ (the bottom of the U-shaped flavor branes), and the instanton's mass is

$$M_I = N_c M_{\text{KK}} \times \left(\lambda B + \sqrt{\frac{2D(1 - C^2)}{5B}} + O(1/\lambda) \right). \quad (2.27)$$

For the antipodal Sakai-Sugimoto model

$$a = \frac{\sqrt[4]{162\pi/5} \approx 3.2}{M_{\text{KK}} \sqrt{\lambda}} \quad \text{and} \quad M_I = N_c M_{\text{KK}} \times \left(\frac{\lambda}{27\pi} + \sqrt{\frac{18\pi}{5}} + O(1/\lambda) \right), \quad (2.28)$$

while other models should have different $O(1)$ numeric factors.

We may absorb 2 out of 3 model-dependent parameters B, C, D into a redefinition of the λ and M_{KK} parameters of the effective 5D theory, for example

$$\lambda \rightarrow \frac{\lambda}{\sqrt{DB^3}}, \quad M_{\text{KK}} \rightarrow M = M_{\text{KK}} \times \sqrt{\frac{D}{B}}, \quad (2.29)$$

thus

$$\frac{8\pi^2}{g^2(z)} = N_c \lambda M \left(1 + M^2 \times z^2 + \dots \right). \quad (2.30)$$

For the *static* instantons or multi-instanton systems we may also get rid of the third model-dependent parameter C . Indeed, for the static systems $\Phi(\mathbf{x}, z) = C \times A_0(\mathbf{x}, z)$, so the only effect of the scalar fields is to reduce the net Coulomb force by a constant factor $(1 - C^2)$. We may simulate this effect without any scalar fields by using different gauge couplings

for the electric and magnetic fields, $g_{\text{el}}^2 = (1 - C^2) \times g_{\text{mag}}^2$, or equivalently by rescaling the time dimension x^0 relative to the space dimensions (x^1, x^2, x^3, z) ,

$$t \rightarrow \frac{t}{\sqrt{1 - C^2}} \quad \text{but} \quad \mathbf{x} \rightarrow \mathbf{x}, \quad z \rightarrow z. \quad (2.31)$$

Consequently, the static instanton's energy becomes

$$\begin{aligned} E_{\text{net}} &= N_c \lambda M + N_c \lambda M^3 \times \int d^3 \mathbf{x} dz I(\mathbf{x}, z) \times z^2 \\ &\quad + \frac{N_c}{4 \lambda M} \times \int d^3 \mathbf{x} dz d^3 \mathbf{x}' dz' \frac{I(\mathbf{x}, z) \times I(\mathbf{x}', z')}{(\mathbf{x} - \mathbf{x}')^2 + (z - z')^2} + \text{subleading} \\ &= N_c \lambda M + N_c \lambda M^3 \times \left(Z^2 + \frac{a^2}{2} \right) + \frac{N_c}{5 \lambda M a^2} + \text{subleading}, \end{aligned} \quad (2.32)$$

hence in equilibrium $Z = 0$ and

$$a = \frac{\sqrt[4]{2/5}}{M \sqrt{\lambda}}. \quad (2.33)$$

Besides the radius a and Z , the instanton has other moduli — the $X^{1,2,3}$ coordinates of the center (which corresponds to the baryon's coordinates in 3D) and $4N_f - 5$ orientation moduli in the $\text{SU}(N_f)$ gauge algebra; and the net energy is degenerate in these moduli to all orders in $1/\lambda$. For finite N_c — even if it's very large — one should quantize the motion of the instanton in those moduli directions. Consequently, a holographic baryon acquires definite spin and isospin quantum numbers; for $N_f = 2$ the baryons have $I = J = 0, 1, 2, \dots, \frac{N_c}{2}$ for even N_c or $I = J = \frac{1}{2}, \frac{3}{2}, \frac{5}{2}, \dots, \frac{N_c}{2}$ for odd N_c [5, 40], and there are similar spin \leftrightarrow flavor multiplet relations for $N_f > 2$.

However, in multi-baryon systems interactions between the baryons break the rotational and flavor symmetries of individual baryons, and only the overall $\text{SO}(3)$ and $\text{SU}(N_f)$ symmetries remain unbroken. In holography, the multi-instanton systems suffer from the same problem: the magnetic fields of multiple instantons interfere with each other, which spoils the degeneracy of the net energy with respect to orientations moduli of the individual instantons. In the large N_c limit, this effect becomes more important than the quantum motion in the moduli space.

Consequently, in our research program into holographic nuclear matter and related issues, we stick to classical static instantons with definite classical orientations in space and in $\text{SU}(N_f)$. From the quantum point of view, such instantons are superpositions of states with different spins and isospins (or rather $\text{SO}(4)$ and $\text{SU}(N_f)$ quantum numbers). We do not even care if the instantons are bosons or fermions — at our level of analysis it simply does not matter. Minimizing the classical energy of a classical multi-instanton system with respect to classical positions and orientations of all the instantons is already a very hard problem — see the next section for the highlights, — and we do not wish to delve into the quantization of moduli until we have the classical issues under control.

2.5 Multi-baryon systems

In the large N_c limit, nuclear forces between the baryons are dominated by the static potentials. Holographically, a static system of A baryons corresponds to a time-independent

configuration of the non-abelian magnetic flavor fields $F_{\mu\nu}^a(\mathbf{x}, z)$ ($\mu, \nu = 1, 2, 3, z$) of net instanton number A ,

$$\int d^3\mathbf{x} dz \frac{\epsilon^{\kappa\lambda\mu\nu}}{16\pi^2} \text{tr}(F_{\kappa\lambda}F_{\mu\nu}) = A = \#\text{baryons}, \quad (2.34)$$

accompanied by the Coulomb electric $A_0^a(\mathbf{x}, z)$ and scalar $\Phi^a(\mathbf{x}, z)$ potentials induced by their Chern-Simons and ΦFF couplings to the magnetic fields [2]. The whole configuration should minimize the net DBI + CS energy of the system subject to the constraint (2.34).

In the $\lambda \rightarrow \infty$ limit, the DBI energy of the magnetic fields is $O(\lambda)$ while the net effect of the electric and scalar fields is only $O(1)$. Moreover, the magnetic fields are concentrated within $O(\lambda^{-1/2})$ distance from the $X^4 = 0$ hyperplane, so to the leading order we may approximate the $4 + 1$ dimensional spacetime as flat. Thus,

$$E_{\text{net}} \approx E_{\text{DBI}} \approx \frac{1}{g_{\text{YM}}^2} \int d^3\mathbf{x} dz \text{Str} \left[\sqrt{\det(K\delta_{\mu\nu} + F_{\mu\nu})} - K \right] \quad (2.35)$$

where we neglect the z dependence of the g_{YM}^2 and $K = g_{11}(z)/(2\pi\alpha')$. Similar to the Yang-Mills energy of an A instanton system, this leading-order DBI energy is minimized by the self-dual configurations of the magnetic fields $F_{\mu\nu}^a(\mathbf{x}, z)$. In fact, all such self-dual configurations (of the same instanton number A) have the same leading-order energies

$$E_{\text{leading order}} = A \times \left(\frac{8\pi^2}{g_{\text{YM}}^2} = N_c \lambda M \right). \quad (2.36)$$

The self-dual configurations form a continuous family parametrized by $A \times 4N_f$ moduli, which correspond to the locations, radii, and $\text{SU}(N_f)$ orientations of the A instantons. But the leading-order energy (2.36) does not depend on any of these moduli.

Fortunately, the sub-leading corrections to the net energy lift the degeneracy of the leading order, which provides for the $O(\lambda^0)$ interactions between the baryons. To work out such interactions we need the degenerate perturbation theory for the magnetic field configurations and their energies. At first order of this perturbation theory, we (a) limit the $F_{\mu\nu}^a(\mathbf{x}, z)$ configurations to the degenerate minima of the leading-order energy function, i.e., to the self-dual magnetic fields; (b) calculate the $O(\lambda^0)$ corrections ΔE for the energies of these configurations; (c) minimize the ΔE among the self-dual configurations. At the next order, we would calculate the $O(\lambda^{-1})$ corrections $\Delta F_{\mu\nu}^a(\mathbf{x}, z)$ to the magnetic fields — which would no longer be self-dual — and then use such corrections to calculate the net energy to the order $O(\lambda^{-1})$. But in this article, we limit our calculations to the $O(\lambda^0)$ interaction energies and zeroth-order magnetic fields, so we shall stop after step (c).

In more detail, to get $O(\lambda^0)$ interactions between A baryons, we proceed as follows:

1. General self-dual magnetic field configurations obtain via ADHM construction [16] in terms of $A \times A$ and $A \times N_f$ matrices obeying certain quadratic constraints. Our first task is to solve these constraints and write down the ADHM matrices in terms of the instantons' locations, radii, and orientations.

2. Given the ADHM matrices, we work out the instanton number density profile

$$I(\mathbf{x}, z) = \frac{\epsilon^{\kappa\lambda\mu\nu}}{16\pi^2} \text{tr}(F_{\kappa\lambda}F_{\mu\nu}), \quad (2.37)$$

and for $N_f > 2$ also the non-abelian adjoint density

$$I^a(\mathbf{x}, x) = \frac{\epsilon^{\kappa\lambda\mu\nu}}{16\pi^2} \times d^{abc} F_{\kappa\lambda}^b F_{\mu\nu}^c. \quad (2.38)$$

3. Next, we calculate the $O(\lambda^0 \times N_c M)$ corrections to the net energy of the system. There are two sources of such corrections: the z -dependence of the 5D gauge coupling, and the Coulomb electric and scalar potentials induced by the Chern-Simons and the ΦFF couplings to the magnetic fields, thus

$$\Delta E = \Delta E_{NA} + \Delta E_C. \quad (2.39)$$

The z -dependent 5D gauge coupling changes the DBI energy of the magnetic fields by

$$\Delta E_{NA} = N_c M \times \int d^3 x dz \lambda M^2 z^2 \times I(\mathbf{x}, z), \quad (2.40)$$

while the Coulomb energy depends on the N_f . For $N_f = 2$, the U(2) Chern-Simons and ΦFF terms couple the SU(2) magnetic fields to the U(1) Coulomb fields only. Consequently, the A_0^a and the Φ^a fields are abelian and couple to the instanton density $I(\mathbf{x}, z)$, and their net energy is simply 4 + 1 dimensional Coulomb energy

$$\Delta E_C = \frac{N_c M}{4} \times \int d^3 \mathbf{x}_1 dx_1 \int d^3 \mathbf{x}_2 dz_2 \frac{I(\mathbf{x}_1, z_1) \times I(\mathbf{x}_2, z_2)}{(\lambda M^2) \times ((\mathbf{x}_1 - \mathbf{x}_2)^2 + (z_1 - z_2)^2)}. \quad (2.41)$$

For $N_f > 2$, the U(N_f) U(2) Chern-Simons and ΦFF terms couple the SU(N_f) magnetic fields to both abelian and non-abelian electric and scalar fields; the abelian Coulomb fields are sourced by the instanton density $I(\mathbf{x}, z)$ while the non-abelian fields are sourced by the adjoint density $I^a(\mathbf{x}, z)$. Altogether, the net energy of these Coulomb fields is

$$\Delta E_C = \frac{N_c M}{4} \times \int d^3 \mathbf{x}_1 dz_1 \int d^3 \mathbf{x}_2 dz_2 \frac{\frac{2}{N_f} I(\mathbf{x}_1, z_1) \times I(\mathbf{x}_2, z_2) + 4 I^a(\mathbf{x}_1, z_1) \times I^a(\mathbf{x}_2, z_2)}{(\lambda M^2) \times ((\mathbf{x}_1 - \mathbf{x}_2)^2 + (z_1 - z_2)^2)}. \quad (2.42)$$

4. Steps 1, 2, and 3, give us ΔE as a function of baryons' locations, radii, and SU(N_f) orientations. Now we need to minimize the ΔE with respect to all these moduli.

This 4-step procedure is fairly straightforward for a few baryons — cf. calculations of the 2-body nuclear forces by Kim, Sin and Zahed [47] and by Hashimoto et al. [48], — but it becomes prohibitively difficult for large numbers of baryons and outright impossible for infinite baryon crystals.⁴ At best, we can survey a small subspace of the A -instanton

⁴A related approach of matrix formulation of multi-D-branes=instantons on flavor branes, was discussed in [49], [50] and [51].

moduli space and try to minimize the ΔE over that subspace. For example, we may assume that all the instantons have the same radius, that their centers form a periodic lattice of some particular symmetry, and that the orientations of the instantons also form some kind of a periodic pattern; this gives us an ansatz for all the $A \times 4N_f$ moduli in terms of just a few overall parameters, and we can try to calculate and minimize the ΔE as a function of these parameters. However, any such ansatz is likely to miss the true lowest-energy configuration of the system. Indeed, in condensed matter guessing the crystalline symmetry of some substance from the properties of the individual atoms is a game of chance with poor odds, and there is no reason why the instanton crystals should be any simpler. Moreover, even if we could somehow guess all the symmetries of the instanton crystal, actually working through steps 1–4 is impossible without additional approximations (besides $\lambda \gg 1$). For example, in our previous paper [1] we needed the (lattice spacing) \gg (instanton radius) approximation just to solve the ADHM constraints for the zigzag-shaped chain of instantons, and even for a straight chain we needed the spacing \gg radius approximation to calculate the Coulomb energy ΔE_C .

In this article we explore a shortcut around steps 1, 2, and 3. In section 3 we shall see that when the distances between the instantons are much larger than their radii, the interactions between the baryons are dominated by the two-body forces,

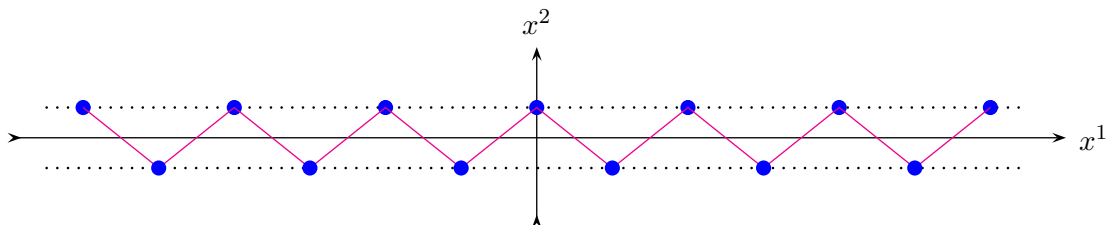
$$\mathcal{E}_{\text{interactions}}^{\text{net}} \approx \frac{1}{2} \sum_{\substack{m,n=1,\dots,A \\ m \neq n}} F_2(X_m^\mu - X_n^\mu, \text{orientation}_m, \text{orientation}_n) \quad (2.43)$$

for a manageably simple function F_2 of the two instanton's positions and orientations. Consequently, we can minimize the net interaction energy over the entire moduli space of the multi-instanton system using a simple numerical simulation: starting from a random set of instanton positions and orientations, use the steepest descent algorithm to find the nearest local minimum of the net energy; repeat this procedure for different random starting points to find other local minima; eventually, we find all the local minima, compare their energies, and identify the global minimum. In our next paper [58] we shall use this algorithm to construct the 2D and 3D instanton lattices starting from scratch, i.e., from the F_2 function. In the present article, we focus on the 1D lattice (section 4) and on the first step in the transition from 1D to 2D — the zigzag (section 5). In both cases, we assume the lattice geometry for the instanton positions, but we use the numeric simulation to find the lowest-energy pattern of their orientations. Once we know all the patterns that show up, we use them as ansatz's (with a few parameters) for which we calculate the net energy as analytic functions of the parameters. In this way, we map the boundaries between different orientation patterns much more accurately than we could do it in the numerical simulation.

To make the baryons form a 1D lattice instead of spreading out in three dimensions, we curve the x_2 and x_3 dimensions of the flavor brane similar to the curvature of the z coordinate. In terms of the effective $4 + 1$ dimensional theory, this corresponds to the 5D flavor gauge coupling depending on the x_2 and the x_3 as well as the $x_4 \equiv z$,

$$\frac{8\pi^2}{g_5^2(x)} = N_c \lambda M \times \left(1 + M_2^2 x_2^2 + M_3^2 x_3^2 + M_4^2 x_4^2 + O(M^4 x^4) \right). \quad (1.2)$$

This gauge coupling acts as a harmonic potential for the instantons which pulls them towards the x_1 axis, so at low densities the instantons form a 1D lattice along the x_1 . At higher densities, the instantons push each other away from the x_1 axis and form more complicated 2D or 3D lattices, starting with the zigzag



To make sure the transition from the straight chain to the zigzag happens for lattice spacings much larger than the instanton radius (which is required by the two-body force approximation), we assume $M_2 \ll M_3, M_4$.

This setting is similar to what we have used in our previous paper [1], except for a change of notations⁵ and allowing for $M_3 \neq M_4$. In section 5 we shall see that $M_3 \neq M_4$ makes a big difference for the instanton orientation patterns in a zigzag: the lowest-energy patterns are abelian for $M_3 \approx M_4$ but non-abelian for anisotropic $M_3 < 0.94M_4$. Even the order of the phase transition from the straight chain to the zigzag depends on the M_3/M_4 ratio: second-order for ratios > 0.725 but first-order for ratios < 0.725 .

3 Two-baryon versus multi-baryon interactions

In real-life nuclear physics, besides the two-body nuclear forces due to meson exchanges, there are significant three-body forces, and presumably also four-body forces, etc.,

$$\hat{H}_{\text{nucleus}} = \sum_{n=1}^A \hat{H}^{1\text{ body}}(n) + \frac{1}{2} \sum_{\substack{\text{different} \\ m,n=1,\dots,A}} \hat{H}^{2\text{ body}}(m,n) + \frac{1}{6} \sum_{\substack{\text{different} \\ \ell,m,n=1,\dots,A}} \hat{H}^{3\text{ body}}(\ell,m,n) + \dots \quad (3.1)$$

where n stands for the quantum numbers of the n^{th} nucleon. Likewise, in the holographic nuclear physics interactions between multiple baryons include two-body forces and also three-body, four-body, etc., forces. Even in the classical infinite-mass limit where the holographic baryons become static instantons of the $SU(N_f)$ gauge fields in 4+1 dimensions, the potential energy (due to $g_{5D} \neq \text{const}$ and due to Chern-Simons interactions) of an A -

⁵Back in [1], we had

$$\frac{8\pi^2}{g_5^2(x)} = N_c \lambda M \times \left(1 + M'^2(x_1^2 + x_2^2) + M^2 x_3^2 + O(M^4 x^4) \right), \quad M' \gg M,$$

so the zigzag was in the (x_3, x_4) plane, with the long axis along the x_4 . Translating between the old notations and those of this paper, we have $(x_1, x_2, x_3, x_4)^{\text{old}} = (x_4, x_3, x_2, x_1)^{\text{new}}$, $M^{\text{old}} = M_2^{\text{new}}$, $M'^{\text{old}} = M_3^{\text{new}} = M_4^{\text{new}}$.

instanton system has form

$$\begin{aligned}\mathcal{E}^{\text{net}}(1, 2, \dots, A) &\equiv E^{\text{total}} - A \times N_c \lambda M \\ &= \sum_{n=1}^A \mathcal{E}^{(1)}(n) + \frac{1}{2} \sum_{\substack{\text{different} \\ m, n=1, \dots, A}} \mathcal{E}^{(2)}(m, n) + \frac{1}{6} \sum_{\substack{\text{different} \\ \ell, m, n=1, \dots, A}} \mathcal{E}^{(3)}(\ell, m, n) + \dots\end{aligned}\quad (3.2)$$

with significant three-body, etc., terms.

What about the relative magnitudes of the 2-body, 3-body, 4-body, etc., interaction terms? When the baryons are packed cheek-by-jowl till their instanton cores overlap and merge, we expect all the n -body forces to have comparable strengths. *But in the opposite low-density regime of baryons separated by distances much larger than their radii, the two-body forces dominate the interactions, while the multi-body forces are smaller by powers of (radius/distance)² [51].* The purpose of this section is to prove this fact.

Before we delve into the technical issues, let us briefly summarize our argument. First, consider the simplest case of $N_f = 2$ and intermediate-range distances between the baryons,

$$\left(a \sim \frac{1}{M\sqrt{\lambda}}\right) \ll |X_m - X_n| \ll \frac{1}{M}, \quad (3.3)$$

which allows us to treat the 5D holographic space as approximately flat. For $N_f = 2$ the holographic baryons are instantons of the SU(2) magnetic fields, which source the U(1) electric and scalar fields via CS and $\text{tr}(\Phi F_{\mu\nu} F^{\mu\nu})$ couplings. When those instantons are small and separated from each other by large distances $|X_m - X_n| \sim D \gg a$, their interactions come from two sources:

1. Direct Coulomb repulsion (electric – scalar) between nearly-point-like abelian charges in $4 + 1$ dimensions,

$$\mathcal{E}^{\text{direct}} = \frac{N_c}{4\lambda M} \sum_{m \neq n} \frac{1}{|X_m^\mu - X_n^\mu|^2}, \quad (3.4)$$

which is a manifestly two-body interaction.

2. The interference between the instantons changes the distribution of the instanton number density in space,

$$I(x) = \frac{\epsilon^{\mu\nu\rho\sigma}}{32\pi^2} \text{tr}(F_{\mu\nu}(x) F_{\rho\sigma}(x)), \quad (3.5)$$

$$I^{A \text{ instantons}}(x) = \sum_{i=n}^A I_n^{\text{standalone}}(x) + \Delta I^{\text{interference}}(x), \quad (3.6)$$

which in turn changes the self-interaction energy (Coulomb and non-abelian) of each instanton by an amount comparable to (3.4). Within an instanton — i.e., at $O(a)$ distance from some instanton's center X_n^μ — the interference from the other instantons should be relatively weak,

$$\Delta I^{\text{interference}}(x) \sim \frac{a^2}{D^2} \times I_n^{\text{standalone}}(x), \quad (3.7)$$

so there should be some kind of a perturbation theory for it. At the first order of such perturbation theory, the ΔI arises from interference between the un-perturbed standalone-like instantons, so we expect it to be a sum of pair-wise interferences from the other instantons,

$$\Delta I^{\text{1st order}}(x \text{ near } X_n) = I_n^{\text{standalone}}(x) \times \sum_{m \neq n} \mathcal{F}_{n,m}^{(1)}(x) \quad (3.8)$$

where

$$\mathcal{F}_{n,m}^{(1)}(x) \sim O\left(\frac{a^2}{|X_n - X_m|^2}\right) \quad (3.9)$$

and depends only on the instantons $\#n$ and $\#m$ — i.e., on their positions, radii, and orientations — but not on any other instantons. At the second-order, we expect to include the interference between the first-order ΔI and the additional instantons, so at this order we obtain 3-body effects,

$$\Delta I^{\text{2nd order}}(x \text{ near } X_n) = I_i^{\text{standalone}}(x) \times \sum_{\ell, m \neq n} \mathcal{F}_{n,\ell,m}^{(2)}(x) \quad (3.10)$$

but

$$\mathcal{F}^{(2)} \sim \frac{a^4}{D^4} \ll \mathcal{F}^{(1)}. \quad (3.11)$$

Likewise, the higher orders may involve more and more instantons, but the magnitudes of such high-order interference effects are suppressed by the higher powers of (a^2/D^2) .

Now consider the Coulomb self-interaction of the instanton $\#n$,

$$\mathcal{E}_C^{\text{self}}(n) = \frac{N_c}{4\lambda M} \times \frac{1}{\rho_n^2} \quad (3.12)$$

where ρ_n is the instanton's effective charge radius,

$$\frac{1}{\rho_n^2} = \iint_{x_1, x_2 \text{ near } X_n} d^4 x_1 d^4 x_2 \frac{I(x_1)I(x_2)}{|x_1 - x_2|^2}. \quad (3.13)$$

A standalone instanton has

$$\frac{1}{\rho_n^2} = \frac{4/5}{a_n^2} \quad (3.14)$$

but interference from the other instantons should change this radius by small amount of similar relative magnitude to $\Delta I/I_n^{\text{standalone}}$, thus

$$\frac{1}{\rho_n^2} \rightarrow \frac{4/5}{a_n^2} + \Delta_n, \quad \Delta_n \sim \frac{1}{D^2}, \quad (3.15)$$

which changes the instanton's Coulomb self-interaction energy by

$$\Delta \mathcal{E}_C^{\text{self}}(n) = \frac{N_c}{4\lambda M} \times \Delta_n \sim \frac{N_c}{\lambda M D^2}. \quad (3.16)$$

Note that this effect has a similar magnitude to the direct Coulomb repulsion (3.4) between the instances.

Moreover, the charge radius correction Δ_n is linear in the $\Delta I^{\text{interference}}$ at x near the X_n , hence in light of eq. (3.8), the leading-order contribution to the Δ_n is a sum of pair-wise interferences from the other instantons, thus

$$\Delta_n = \sum_{m \neq n} \Delta_{n,m}^{(1)} + O(a^2/D^4) \quad (3.17)$$

where each $\Delta_{n,m}^{(1)}$ depends only on the instantons $\#n$ and $\#m$. Consequently, the leading effect of the interference on the net Coulomb self-energy of all the instantons has form

$$\begin{aligned} \Delta^{\text{interference}} \mathcal{E}_C^{\text{self}} &\equiv \mathcal{E}_C^{\text{self}}[\text{interfering}] - \mathcal{E}_C^{\text{self}}[\text{standalone}] \\ &= \frac{N_c}{4\lambda M} \sum_{\substack{n,m=1,\dots,A \\ n \neq m}} \Delta_{n,m}^{(1)} + O\left(\frac{N_c}{\lambda M} \times \frac{a^2}{D^4}\right) \end{aligned} \quad (3.18)$$

where the leading terms act as two-body interactions between the instantons.

Besides the Coulomb energy, the non-abelian energy is also affected by the interference between the instantons,

$$\Delta^{\text{interference}} \mathcal{E}_{NA} = N_c \lambda M^3 \times \int d^4x (x^4)^2 \times \Delta I^{\text{interference}}(x), \quad (3.19)$$

but this time the integral should be taken over the whole 4D space, including both the instantons and the inter-instanton space. Indeed, we shall see later in this section that

$$\begin{aligned} \text{at } O(a) \text{ distance from an } X_n, \quad \Delta I &\sim \frac{1}{a^2 D^2}, \\ \text{at } O(D) \text{ distances from all the } X_m, \quad \Delta I &\sim \frac{a^4}{D^8}, \end{aligned} \quad (3.20)$$

and both kinds of places make $O(a^4/D^2)$ contributions to the integral (3.19). Moreover, in both kinds of places, the leading terms in the $\Delta I(x)$ is a sum of independent two-instanton interference terms,

$$\Delta I^{\text{interference}}(x) = \frac{1}{2} \sum_{n \neq m} \mathcal{I}_{n,m}^{(2)}(x) + \text{subleading}. \quad (3.21)$$

Although our heuristic argument for such decomposition near instanton centers does not work in the inter-instanton space, we shall prove later in this section that the decomposition (3.21) works everywhere in the 4D space. Therefore, the non-abelian interactions between the instantons due to interference are also dominated by the two-body terms

$$\Delta^{\text{interference}} \mathcal{E}_{NA} = \frac{1}{2} \sum_{n \neq m} N_c \lambda M^3 \int d^4x (x^4)^2 \times \mathcal{I}_{n,m}(x) + \text{subleading}. \quad (3.22)$$

Thus far we have assumed $N_f = 2$. For larger numbers of flavors, interference between magnetic fields of the $SU(N_f)$ instantons is harder to describe mathematically, but its qualitative features for $a \ll D$ remain exactly as in eqs. (3.5) through (3.21). In particular, the effect of interference between small instantons separated by large distances on the

density profile $I(x)$ is dominated by the two-instanton interference terms as in eqs. (3.21) and (3.8)–(3.11). Consequently, the instantons' non-abelian and Coulomb interaction energies suffer small corrections from the interference that give rise to predominantly-two-body interactions between the instantons as in eqs. (3.18) and (3.22).

The qualitatively new aspect of $N_f \geq 3$ comes not from the $SU(N_f)$ self-dual magnetic fields of the multiple instantons but from the electric and scalar Coulomb fields induced by the magnetic fields via Chern-Simons and $\text{tr}(\Phi F^{lmuv} F_{\mu\nu})$ interactions. For $N_f = 2$ such electric and scalar fields are purely abelian, $A^0, \Phi \in U(1) \subset U(2)$, but for $N_f \geq 3$ they have both $U(1)$ and $SU(N_f)$ components. For a single instanton, all fields — magnetic, electric, and scalar — belong to the same Cartesian $U(2)$ subgroup of the $U(N_f)$ — in $N_f \times N_f$ matrix language,

$$\text{for } A_{\text{magnetic}}^\mu \in \left(\begin{array}{c|c} SU(2) & \mathbf{0} \\ \hline \mathbf{0} & \mathbf{0} \end{array} \right), \quad A^0, \Phi \propto \begin{pmatrix} \frac{1}{2} & & & \mathbf{0} \\ & \frac{1}{2} & & \\ & & 0 & \\ \mathbf{0} & & & 0 \end{pmatrix} \quad (3.23)$$

— so the Coulomb *self-interaction* of a standalone instanton works exactly as for $N_f = 2$. On the other hand, Coulomb interactions *between* multiple instantons whose magnetic fields span different $SU(2)$ subgroups of the $SU(N_f)$ become much more complicated.

Moreover, in a general multi-instanton background, the Coulomb fields $A^0(x)$ and $\Phi(x)$ do not behave as simply $1/r^2$. Instead, they satisfy gauge-covariant equations for the matrix-valued fields,

$$-\left(\partial_\mu + i[A_\mu(x), \cdot]\right)^2 A^0(x) = \frac{g^2 N_c}{32\pi^2} \epsilon_{\alpha\beta\mu\nu} F^{\alpha\beta} F^{\mu\nu} \quad (3.24)$$

and ditto for the scalar $\Phi(x)$. Consequently, the Coulomb force between two sources (such as instantons) depends on the magnetic fields in the space between them. In other words, for $N_f \geq 3$ the direct Coulomb interactions are no longer two-body; instead, the force between two instantons depends on the other instantons in the system.

Fortunately, for small instantons separated by large distances, the magnetic fields are concentrated in small volumes of 4-space, so their effect on the Coulomb field propagation becomes perturbatively weak:

$$\begin{aligned} \left\langle x^\mu, b \left| \frac{-1}{\text{covariant } \square} \right| y^\mu, c \right\rangle &= \frac{\delta^{bc}}{4\pi^2 |x^\mu - y^\mu|^2} + \sum_{n=1}^A \frac{O(a_n^2) F(b, c, \text{orientation}_n)}{|x^\mu - X_n^\mu|^2 \times |y^\mu - X_n^\mu|^2} \\ &+ O(a^4/D^6) \text{ two-instanton terms} + \dots \end{aligned} \quad (3.25)$$

where b and c are adjoint indices of the $U(N_f)$. The leading term here is the ordinary Coulomb propagator in 4D space, thus *in the $a \ll D$ limit, the direct Coulomb interactions between small instantons are predominantly two-body*,

$$\mathcal{E}^{\text{direct}} \approx \frac{N_c}{4\lambda M} \sum_{n \neq m} \frac{\text{overlap}(n, m)}{|X_n^\mu - X_m^\mu|^2} \quad (3.26)$$

where $\text{overlap}(n, m) \geq 0$ is the overlap between the $\text{SU}(2)$'s spanned by the magnetic fields of the instantons $\#n$ and $\#m$.

The bottom line is, *for all $N_f \geq 2$, all the interactions between small instantons separated by large distances are predominantly two-body*. However, for $N_f \geq 3$ the dependence of such two-body interactions on the instantons' orientations becomes much more complicated than for $N_f = 2$.

Finally, let us relax our second un-necessary assumption of intermediate-range distances between the baryons and consider what happens at longer distances $D \sim 1/M$. In this regime, the curvature of the fourth space dimension — and especially the x^4 -dependence of the 5D gauge coupling — can no longer be treated as a perturbation. Consequently, the magnetic fields of an instanton or a multi-instanton system are no longer self dual in the inter-instanton space — although inside the instanton cores or at intermediate-range distances $r \ll 1/M$ from instantons they remain approximately self-dual. Therefore, the interference between very distant instantons is no longer governed by the self-dual ADHM solutions. Instead, we must work it out the hard way: figure out how the magnetic fields of an instanton propagate through the curved $4 + 1$ dimensions towards the other instantons, and then find out how such fields disturb the other instantons' cores.

Fortunately, we do not need a hard calculation to see that at large distances from an instanton its magnetic fields are very weak. Indeed, even in flat 5D space the fields weaken with distance as $A_\mu^a \sim a^2/r^3$ (in the IR-safe singular gauge), so at $r \gg a$ they are so weak that the field equations become effectively linear. In the curved space, we may decompose the weak 5D gauge fields into 4D mesonic fields, hence at large distances $r \gtrsim 1/M$ from an instanton, its fields become Yukawa-like

$$A_\mu^a(r, Z) \sim \frac{a^2}{r^3} \times \sum_k \Psi_k(z) \Psi_k(z_{\text{inst}}) e^{-m_k r} \ll \frac{a^2}{r^3}. \quad (3.27)$$

Consequently, at the location X_n^μ of any particular instanton, the background fields from the *other* instantons are very weak, and their effect on the instanton $\#n$ itself can be adequately accounted by the first-order perturbation theory. In other words, the effects of other instantons $\#m \neq n$ on the instanton $\#n$ are weak and add up linearly! For the self-interaction energy of the n^{th} instanton, this means

$$\mathcal{E}(n) \approx \mathcal{E}(\text{standalone}) + \sum_{m \neq n} \Delta_m \mathcal{E}(n) \quad (3.28)$$

where the second term gives rise to two-body interaction energies

$$\mathcal{E}_{\text{interference}}^{(2)}(n, m) = \Delta_m \mathcal{E}(n) + \Delta_n \mathcal{E}(m). \quad (3.29)$$

We expect the two-body terms to be rather small — in addition to the usual 5D $1/|X_m - X_n|^2$ factors they should carry Yukawa-like exponentials e^{-mr} (or rather sums of such exponentials), but these are the leading interactions due to interference. To three-body or multi-body interactions follow from higher-order perturbation by very weak fields (3.27), so they are much smaller than the two-body interactions.

As to the direct Coulomb interactions between the holographic baryons via electric or scalar fields, at large distances $|X_m - X_n| \sim 1/M$ they also decompose into sums of Yukawa forces. Moreover, since the scalar mesons generally have different masses from the vector mesons, the attractive potential due to scalars may have a different r dependence from the repulsive potential due to vectors. Thus, for a right model, the net two-body force between two holographic baryons may become attractive at large enough distances between them. But regardless of the model, the direct Coulomb interactions are always manifestly two-body for $N_f = 2$, while for $N_f \geq 3$ the multi-body terms exist but become very small at large distances between the baryons. Thus,

$$\mathcal{E}_{\text{direct}} = \frac{1}{2} \sum_{m \neq n} \text{overlap}(m, n) \times V^{(2)}(|X_m - X_n|) \quad (3.30)$$

where the precise form of the potential $V^{(2)}(r)$ is model-dependent, but the two-body form of the direct interactions is quite universal.

★ ★ ★

At this point we have finished our heuristic arguments and about to start a more technical analysis of the interactions between the holographic baryons. To keep our presentation relatively simple, we assume just two quark flavors and intermediate-range distances between the holographic baryons — $D \gg a$ but $D \ll 1/M$.

For $N_f = 2$ the non-abelian gauge symmetry is $SU(2)$, so the ADHM data for an A -instanton configuration consists of a quaternionic array Y_n ($n = 1, \dots, A$) and a symmetric quaternionic $A \times A$ matrix Γ_{mn} . Equivalently, we may use A real instanton radii $a_n = |Y_n|$, A $SU(2)$ matrices y_n (equivalent to unimodular quaternions Y_n/a_n) parametrizing the instanton's orientations, and 4 real symmetric matrices Γ_{mn}^μ whose diagonal elements $\Gamma_{nn}^\mu = X_n^\mu$ are the locations of the instanton's centers. Finally, the off-diagonal matrix elements $\Gamma_{m \neq n}^\mu = \alpha_{mn}^\mu$ follow from the other parameters by solving the ADHM equations

$$\Im \left((\Gamma^\dagger \Gamma)_{mn} + Y_m^\dagger Y_n \right) = 0 \quad (3.31)$$

or equivalently

$$\eta_{\mu\nu}^a [\Gamma^\mu, \Gamma^\nu]_{mn} + a_m a_n \times \text{tr}(y_m^\dagger y_n (-i\tau^a)) = 0 \quad (3.32)$$

where $\eta_{\mu\nu}^a$ is the 't Hooft's tensor mapping between the $SU(2)_{\text{gauge}}$ and the $SU(2)_L$ inside $\text{Spin}(4) = SU(2)_L \times SU(2)_R$,

$$\eta_{44}^a = 0, \quad \eta_{4i}^a = -\delta_i^a, \quad \eta_{i4}^a = +\delta_i^a, \quad \eta_{ij}^a = \epsilon^{aij}, \quad a, i, j = 1, 2, 3. \quad (3.33)$$

The ADHM data are somewhat redundant — an $O(A)$ symmetry acting on all the Γ_{mn}^μ and $Y_n = a_n y_n$ does not change any physical properties of the multi-instanton data. This symmetry includes \mathbb{Z}_2^A which flips the instanton orientations $y_n \rightarrow -y_n$ (independently for each n). It also includes small $SO(A)$ rotations that change the off-diagonal elements by $\delta\alpha_{mn}^\mu = \epsilon_{mn}(X_m - X_n)^\mu + O(\epsilon^2)$. To eliminate these rotations, the ADHM equations (3.32)

for the off-diagonal elements should be combined with additional constraints (one for each $m \neq n$), for example

$$\forall m \neq n : (X_m - X_n)^\mu \alpha_{mn}^\mu = 0. \quad (3.34)$$

For large distances $|X_m - X_n| \sim D$ between the instantons, $D \gg a_n$, we may solve the ADHM equations (and the constraints (3.34) as a power series in a^2/D^2 :

$$\alpha_{mn}^\mu \equiv \Gamma_{m \neq n}^\mu = \alpha_{mn}^{(1)} + \alpha_{mn}^{(2)} + \alpha_{mn}^{(3)} + \dots, \quad (3.35a)$$

$$\alpha_{mn}^{(1)} = \frac{\eta_{\mu\nu}^a (X_m - X_n)_\nu}{|X_m - X_n|^2} \times \frac{1}{2} a_m a_n \text{tr}(y_m^\dagger y_n (-i\tau^a)) = O(a^2/D), \quad (3.35b)$$

$$\alpha_{mn}^{(2)} = -\frac{\eta_{\mu\nu}^a (X_m - X_n)_\nu}{|X_m - X_n|^2} \times \sum_{\ell \neq m, n} \eta_{\kappa\lambda}^a \alpha_{\ell m}^{(1)} \alpha_{\ell n}^{(1)} = O(a^4/D^3), \quad (3.35c)$$

$$\alpha_{mn}^{(3)} = -2 \frac{\eta_{\mu\nu}^a (X_m - X_n)_\nu}{|X_m - X_n|^2} \times \sum_{\ell \neq m, n} \eta_{\kappa\lambda}^a \alpha_{\ell m}^{(1)} \alpha_{\ell n}^{(2)} = O(a^6/D^5), \quad (3.35d)$$

...

Note that for each off-diagonal matrix element, the leading term $\alpha_{mn}^{(1)}$ in this expansion depends only on the instantons $\#m$ and $\#n$ (i.e., on the positions, radii, and orientations of only these two instantons) while the subleading terms $\alpha_{mn}^{(2)}, \alpha_{mn}^{(3)}, \dots$ involve additional instantons.

Given the ADHM data Γ_{mn}^μ , a_n , and y_n of a multi-instanton system, the 4D instanton number density (3.5) can be obtained as

$$I(x) = -\frac{1}{16\pi^2} \square\square \log \det(L(x)) \quad (3.36)$$

where $L(x)$ is a real $A \times A$ symmetric matrix

$$L_{mn}(x) = \sum_{\ell} (\Gamma_{\ell m}^\mu - x^\mu \delta_{\ell m}) (\Gamma_{\ell n}^\mu - x^\mu \delta_{\ell n}) + \frac{1}{2} a_m a_n \text{tr}(y_m^\dagger y_n). \quad (3.37)$$

Thanks to the double D'Alembertian in eq. (3.36), several moments of the instanton density may be obtained via integrating by parts:

$$\int d^4x I(x) = A, \quad (3.38a)$$

$$\int d^4x I(x) \times x^\nu = \text{tr}(\Gamma^\nu), \quad (3.38b)$$

$$\int d^4x I(x) \times x^\mu x^\nu = \text{tr}(\Gamma^\mu \Gamma^\nu) + \frac{1}{2} \delta^{\mu\nu} \text{tr}(T) \quad (3.38c)$$

$$\text{where } T_{mn} \equiv \frac{1}{2} a_m a_n \text{tr}(y_m^\dagger y_n), \quad (3.38d)$$

$$\begin{aligned} \int d^4x I(x) \times x^\lambda x^\mu x^\nu &= \frac{1}{2} (\Gamma^\lambda \{\Gamma^\mu, \Gamma^\nu\}) + \frac{1}{2} \delta^{\lambda\mu} \text{tr}(\Gamma^\nu T) \\ &\quad + \frac{1}{2} \delta^{\lambda\nu} \text{tr}(\Gamma^\mu T) + \frac{1}{2} \delta^{\mu\nu} \text{tr}(\Gamma^\lambda T). \end{aligned} \quad (3.38e)$$

The quadratic moment (3.38c) for $\mu = \nu = 4$ is particularly important since it gives us an exact formula for the non-abelian energy of the system directly in terms of the ADHM data,

$$\begin{aligned}
 \mathcal{E}_{NA} &= N_c \lambda M^3 \times \int d^4 x I(x) \times (x^4)^2 = N_c \lambda M^3 \times \left(\text{tr}(\Gamma^4 \Gamma^4) + \frac{1}{2} \text{tr}(T) \right) \\
 &= N_c \lambda M^3 \times \left(\sum_{i=n}^A \left((\Gamma_{nn}^4)^2 + \frac{1}{2} T_{nn} \right) + \sum_{m \neq n} (\Gamma_{mn}^4)^2 \right) \\
 &= N_c \lambda M^3 \sum_n \left((X_n^4)^2 + \frac{1}{2} a_n^2 \right) + N_c \lambda M^3 \sum_{m \neq n} (\alpha_{mn}^4)^2.
 \end{aligned} \tag{3.39}$$

Obviously, the first sum on the last line here is the sum of individual instantons' potential energies due to their radii and locations (relative to the $x^4 = 0$ hyperplane) while the second sum comprises the interactions between the instantons,

$$\mathcal{E}_{NA}^{\text{net interaction}} = N_c \lambda M^3 \sum_{m \neq n} (\alpha_{mn}^4)^2. \tag{3.40}$$

This formula is general and applies to any multi-instanton configuration, the problem is calculating the off-diagonal matrix elements $\alpha_{mn}^\mu = \Gamma_{m \neq n}^\mu$. Fortunately, for the small instantons separated by large distances $O(D) \gg a$, those off-diagonal matrix elements are given by the expansion (3.35.a-d) in powers of (a^2/D^2) . Moreover, for each $m \neq n$ the leading term $\alpha_{\mu mn}^{(1)}$ in this expansion depends only on the positions and orientations of the instantons $\#m$ and $\#n$ but does not depend on any other instantons. Thus, *to the leading order in a^2/D^2 , the non-abelian interactions (3.40) between the instantons are two-body*, i.e. involve only two instantons at a time:

$$\mathcal{E}_{NA}^{\text{interaction}} \approx \frac{1}{2} \sum_{m \neq n} \mathcal{E}_{NA}^{2\text{body}}(m, n) \sim N_c \lambda M^3 \frac{a^4}{D^2} \tag{3.41}$$

while the multi-body interactions are $O(N_c \lambda M^3 a^6/D^4)$ or weaker. Specifically,

$$\begin{aligned}
 \mathcal{E}_{NA}^{2\text{body}}(m, n) &= 2N_c \lambda M^3 \times \left(a_m a_n \times \frac{\eta_{4\nu}^a (X_m - X_n)_\nu}{|X_m - X_n|_{4D}^2} \times \frac{1}{2} \text{tr}(y_m^\dagger y_n (-i\tau^a)) \right)^2 \\
 &= \frac{N_c \lambda M^3 a_m^2 a_n^2}{2|X_m - X_n|_{4D}^2} \times \text{tr}^2 \left(y_m^\dagger y_n (-i\vec{N}_{mn} \cdot \vec{\tau})_{3D} \right)
 \end{aligned} \tag{3.42}$$

where \vec{N}_{mn} is the 3-vector part of the unit 4-vector

$$N_{mn}^\mu \equiv (\vec{N}_{mn}, N_{mn}^4) = \frac{X_n^\mu - X_m^\mu}{|X_n - X_m|}. \tag{3.43}$$

Now consider the Coulomb energy

$$\mathcal{E}_C = \frac{N_c}{4\lambda M} \iint d^4 x_1 d^4 x_2 \frac{I(x_1) I(x_2)}{|x_1 - x_2|^2}. \tag{3.44}$$

of the multi-instanton system. Alas, we cannot obtain this Coulomb energy directly from the ADHM data via integration by parts or some other clever trick but have to calculate

it the hard way: derive the instanton density profile from eq. (3.36), decompose $I(x)$ into individual instantons plus interference terms, and finally perform the integral (3.44) for some approximation to the integrand.

Our starting point is the $L_{mn}(x)$ matrix (3.37) which governs the instanton density profile. For a system of small instantons separated by large distances, the diagonal elements of this matrix are much larger than the off-diagonal elements. Indeed, the diagonal matrix elements are given by

$$L_{nn}(x) = \left((\Gamma^\mu - x^\mu)^2 \right)_{nn} + T_{nn} = (X_n^\mu - x^\mu)^2 + \sum_{\ell \neq n} |\alpha_{\ell n}^\mu|^2 + a_n^2 \quad (3.45)$$

and their magnitudes are generally $O(D^2)$. To be precise, in the inter-instanton space all of the $L_{nn}(x)$ are $O(D^2)$ or larger, while in the neighborhood of an instanton $\#m$ the corresponding $L_{mm}(x)$ may become as small as $O(a^2)$ but all the other $L_{nn}(x)$ remain $O(D^2)$. On the other hand, the off-diagonal matrix elements are only $O(a^2) \ll D^2$,

$$\begin{aligned} L_{m \neq n} &= \left((\Gamma^\mu - x^\mu)^2 \right)_{mn} + T_{mn} \\ &= (X_m + X_n - 2x)^\mu \alpha_{mn}^\mu + \sum_{\ell \neq m, n} \alpha_{\ell m}^\mu \alpha_{\ell n}^\mu + a_m a_n \times \frac{1}{2} \text{tr}(y_m^\dagger y_n) \\ &= O(a^2). \end{aligned} \quad (3.46)$$

This hierarchy between the diagonal and the off-diagonal matrix elements of L allows us to evaluate $\log \det(L)$ as a power series in the ratios of the off-diagonal to diagonal elements:

$$\begin{aligned} \log \det(L) &= \log \det(L_{\text{diag}}) + \log \det(1 + L_{\text{diag}}^{-1} L_{\text{offdiag}}) \\ &= \text{tr} \log(L_{\text{diag}}) + \text{tr} \log(1 + L_{\text{diag}}^{-1} L_{\text{offdiag}}) \\ &= \text{tr} \log(L_{\text{diag}}) + \text{tr}(L_{\text{diag}}^{-1} L_{\text{offdiag}}) - \frac{1}{2} \text{tr}(L_{\text{diag}}^{-1} L_{\text{offdiag}} L_{\text{diag}}^{-1} L_{\text{offdiag}}) \\ &\quad + \frac{1}{3} \text{tr}(L_{\text{diag}}^{-1} L_{\text{offdiag}} L_{\text{diag}}^{-1} L_{\text{offdiag}} L_{\text{diag}}^{-1} L_{\text{offdiag}}) + \dots \\ &= \sum_n \log(L_{nn}) + 0 - \frac{1}{2} \sum_{m \neq n} \frac{L_{mn} L_{nm}}{L_{mm} L_{nn}} + \frac{1}{3} \sum_{\substack{\text{different} \\ \ell, m, n}} \frac{L_{\ell m} L_{mn} L_{n \ell}}{L_{\ell \ell} L_{mm} L_{nn}} + \dots \end{aligned} \quad (3.47)$$

Consequently, the net instanton density profile of the system may be written as a sum of one-instanton, two-instanton, three-instanton, etc., terms:

$$\begin{aligned} I(x) &= \frac{-1}{16\pi^2} \square \square \log \det(L) \\ &= \sum_n \mathcal{I}_n^{(1)}(x) + \frac{1}{2} \sum_{m \neq n} \mathcal{I}_{mn}^{(2)}(x) + \frac{1}{6} \sum_{\substack{\text{different} \\ \ell, m, n}} \mathcal{I}_{\ell mn}^{(3)}(x) + \dots \end{aligned} \quad (3.48)$$

where

$$\mathcal{I}_n^{(1)}(x) = \frac{-1}{16\pi^2} \square\square \log(L_{nn}(x)), \quad (3.49)$$

$$\mathcal{I}_{mn}^{(2)}(x) = \frac{+1}{16\pi^2} \square\square \frac{L_{mn}L_{nm}}{L_{mm}L_{nn}}, \quad (3.50)$$

$$\mathcal{I}_{\ell mn}^{(3)}(x) = \frac{-2}{16\pi^2} \square\square \frac{L_{\ell m}L_{mn}L_{n\ell}}{L_{\ell\ell}L_{mm}L_{nn}}, \quad (3.51)$$

etc., etc.

The one-instanton terms (3.49) here have the form of stand-alone instanton profiles. Indeed, rewriting eq. (3.45) as

$$L_{nn}(x) = r_n^2 + \hat{a}_n^2 \quad (3.52)$$

$$\text{where } r_n^2 \equiv |x^\mu - X_n^\mu|^2 \quad \text{and} \quad \hat{a}_n^2 = a_n^2 + \sum_{m \neq n} |\alpha_{mn}^\mu|^2 = a_n^2 + O(a^4/D^2),$$

we obtain

$$\frac{-1}{16\pi^2} \square\square \log(L_{nn}) = \frac{-1}{16\pi^2} \left(\frac{\partial^2}{\partial r_n^2} + \frac{3}{r_n} \frac{\partial}{\partial r_n} \right)^2 \log(r_n^2 + \hat{a}_n^2) = \frac{(6/\pi^2) \hat{a}_n^4}{(r_n^2 + \hat{a}_n^2)^4}, \quad (3.53)$$

which is precisely the density profile of a standalone instanton of radius \hat{a}_n . Note that this radius is slightly larger than the original instanton radius a_n — this is a hidden effect of the interference from the other instantons $m \neq n$ — while the visible effects of the interference come via the two-instanton terms $\mathcal{I}_{mn}^{(2)}(x)$, the three-instanton terms $\mathcal{I}_{\ell mn}^{(3)}(x)$, etc. Deriving explicit formulae for all these terms is a straightforward (albeit rather tedious) exercise, and the results are too cumbersome to present here. Instead, let us simply estimate the magnitudes of the multi-instanton terms.

In the inter-instanton space — i.e. at $O(D)$ distances from *all* the instanton centers — we have

$$\frac{L_{mn}L_{nm}}{L_{mm}L_{nn}} \sim \frac{a^4}{D^4}, \quad \frac{L_{\ell m}L_{mn}L_{n\ell}}{L_{\ell\ell}L_{mm}L_{nn}} \sim \frac{a^6}{D^6}, \quad \text{etc.} \quad (3.54)$$

Moreover, all these expressions change rather slowly with x — on the scale $x \sim D$ — so in the context of such expressions $\partial_\mu = O(1/D)$. Consequently,

$$\square\square \frac{L_{mn}L_{nm}}{L_{mm}L_{nn}} \sim \frac{a^4}{D^8}, \quad \square\square \frac{L_{\ell m}L_{mn}L_{n\ell}}{L_{\ell\ell}L_{mm}L_{nn}} \sim \frac{a^6}{D^{10}}, \quad \dots, \quad (3.55)$$

and therefore

$$\text{In the inter-instanton space} \quad \mathcal{I}_{mn}^{(2)}(x) \sim \frac{a^4}{D^8}, \quad \mathcal{I}_{\ell mn}^{(3)}(x) \sim \frac{a^6}{D^{10}}, \quad \dots \quad (3.56)$$

As promised in eq. (3.21), the interference between small well-separated instantons is dominated by the two-instanton terms. In fact, in the inter-instanton space the two-instanton terms have similar magnitudes to the one-instanton terms,

$$\mathcal{I}_{mn}^{(2)}(x) \sim \frac{a^4}{D^8} \sim \mathcal{I}_n^{(1)}(x). \quad (3.57)$$

However, the entire $I(x)$ in the inter-instanton space makes a negligible contribution to the net Coulomb energy of the system,

$$\mathcal{E}_C[\text{IIS}] = \int_{\text{IIS}} d^4x I(x) \times \frac{N_c}{2\lambda M} \sum_n \frac{1}{|x^\mu - X_n^\mu|^2} \sim D^4 \times \frac{a^4}{D^8} \times \frac{N_c}{\lambda M D^2} \sim \frac{N_c a^4}{\lambda M D^6}. \quad (3.58)$$

Therefore, the net Coulomb energy of the multi-instanton system may be summarized as

$$\mathcal{E}_C^{\text{net}} = \frac{N_c}{4\lambda M} \times \left\{ \sum_n \frac{1}{\rho_n^2} + \sum_{m \neq n} \frac{1}{|X_m^\mu - X_n^\mu|^2} + O(a^4/D^6) \right\}, \quad (3.59)$$

where the first term inside the ‘ $\{\dots\}$ ’ comprises the self-interactions of individual instantons — but with the effective charge radii ρ_n that include the effects of interference from the other instantons, — the second term is the direct Coulomb repulsion (3.4) between different instantons approximated as point charges (or equivalently, as compact spherically-symmetric charges), and the third term accounts for the $O(a^4/D^8)$ instanton density in the inter-instanton space. Note that the self-interaction terms dominate the net Coulomb energy, so even relatively small corrections to the effective charge radii ρ_i can have an effect comparable to the direct Coulomb repulsion between the instantons.

To work out the effects of interference on the charge radius ρ_n of the n^{th} instanton we need to estimate (and then evaluate) the interference terms $\mathcal{I}_{\ell m}^{(2)}$, $\mathcal{I}_{k\ell m}^{(3)}$, etc., for x^μ near the n^{th} instanton’s center X_n^μ . In this neighborhood, the estimates (3.56) do not work — or rather they work only for the interference terms between the *other instantons* $\#k, \ell, m, \dots \neq n$.

But the terms $\mathcal{I}_{mn}^{(2)}$, $\mathcal{I}_{\ell mn}^{(3)}$, etc., that do involve the n^{th} instanton itself turn out to be much larger than (3.56) because of the $1/L_{nn}(x)$ factor that happens to be $O(1/a^2)$ instead of $O(1/D^2)$. Moreover, this factor depends on x much more rapidly than all the other factors — on the scale of $x \sim a$ instead of $x \sim D$ — so taking the space derivatives of this factor produces much larger results than taking the same derivatives of the other factors. Thus

$$\begin{aligned} \square\square \frac{L_{mn}L_{nm}}{L_{mm}L_{nn}} &\approx \left(\square\square \frac{1}{L_{nn}} \right) \times \frac{L_{mn}L_{nm}}{L_{mm}}, \\ \square\square \frac{L_{\ell m}L_{mn}L_{n\ell}}{L_{\ell\ell}L_{mm}L_{nn}} &\approx \left(\square\square \frac{1}{L_{nn}} \right) \times \frac{L_{\ell m}L_{mn}L_{n\ell}}{L_{\ell\ell}L_{mm}}, \\ &\text{etc.}, \end{aligned} \quad (3.60)$$

and therefore

$$\mathcal{I}_{mn}^{(2)}(x) = \left(\frac{1}{16\pi^2} \square\square \frac{1}{L_{nn}} \right) \times \left[\frac{L_{mn}L_{nm}}{L_{mm}} \right] (@x = X_n) + \text{subleading}, \quad (3.61)$$

$$\mathcal{I}_{\ell mn}^{(3)}(x) = \left(\frac{1}{16\pi^2} \square\square \frac{1}{L_{nn}} \right) \times \left[-2 \frac{L_{\ell m}L_{mn}L_{n\ell}}{L_{\ell\ell}L_{mm}} \right] (@x = X_i) + \text{subleading}, \quad (3.62)$$

etc., etc.

The first factor in all these interference terms evaluates to

$$\frac{1}{16\pi^2} \square\square \left(\frac{1}{L_{nn}} = \frac{1}{r_n^2 + \hat{a}_n^2} \right) = \frac{12}{\pi^2} \frac{\hat{a}_n^2(\hat{a}_n^2 + r_n^2)}{(\hat{a}_n^2 + r_n^2)^5} \sim \frac{1}{a^6} \quad (3.63)$$

hence

$$\text{In the vicinity of the } n^{\text{th}} \text{ instanton} \quad \mathcal{I}_{mn}^{(2)}(x) \sim \frac{1}{a^2 D^2}, \quad \mathcal{I}_{\ell mn}^{(3)}(x) \sim \frac{1}{D^4}, \quad \dots \quad (3.64)$$

Since the standalone instanton's density is $I_n^{(1)}(x) \sim 1/a^4$, the interference terms change the effective charge radius or rather

$$\frac{1}{\rho_n} [\text{due to } I_n^{(1)} \text{ only}] = \iint_{\text{near } X_n} d^4x d^4x' \frac{I_n^{(1)}(x) \times I_n^{(1)}(x')}{|x - x'|^2} = \frac{4/5}{\hat{a}_n^2} \quad (3.65)$$

by

$$\Delta_n [\text{due to } \mathcal{I}_{mn}^2] = \iint_{\text{near } X_n} d^4x d^4x' \frac{2I_n^{(1)}(x) \times I_{mn}^{(2)}(x')}{|x - x'|^2} \sim \frac{a^8}{a^2} \times \frac{1}{a^4} \times \frac{1}{a^2 D^2} \sim \frac{1}{D^2}, \quad (3.66)$$

$$\Delta_n [\text{due to } \mathcal{I}_{\ell mn}^3] = \iint_{\text{near } X_n} d^4x d^4x' \frac{2I_n^{(1)}(x) \times I_{\ell mn}^{(3)}(x')}{|x - x'|^2} \sim \frac{a^8}{a^2} \times \frac{1}{a^4} \times \frac{1}{D^4} \sim \frac{a^2}{D^4}, \quad (3.67)$$

etc., etc.

Therefore, *the two-instanton interference terms \mathcal{I}_{mn}^2 affect the Coulomb energy by the amount comparable to the direct repulsion (between point-like instantons) but the effect of interference terms involving three or more instantons at once is suppressed by extra powers of a^2/D^2 and becomes negligible in the limit of small instantons separated by large distances.* Since we have already verified similar behavior of the non-abelian energy of the multi-instanton system, this completes the proof of our theorem. *Quod erat demonstrandum.*

To conclude this section, we would like to derive explicit formulae for the two-instanton interference terms (near one of the two instantons) and hence the Coulomb interaction energy due to interference. Starting from eq. (3.61) we have the first factor on the r.h.s. evaluated in eq. (3.63), so let us evaluate the second factor. For $x^\mu = X_m^\mu$, eq. (3.46) gives us

$$L_{mn}(x = X_I) = L_{nm}(x = X_I) = (X_m - X_n)^\mu \alpha_{mn}^\mu + \sum_{\ell \neq m, n} \alpha_{\ell m}^\mu \alpha_{\ell n}^\mu + a_m a_n \times \frac{1}{2} \text{tr}(y_m^\dagger y_n) \quad (3.68)$$

where the first term on the r.h.s. happens to vanish — cf. eqs. (3.34) — while the second term is $O(a^4/D^2)$, which is much smaller than the $O(a^2)$ third term. Thus,

$$L_{mn}(x = X_n) = L_{nm}(x = X_n) \approx a_m a_n \times \frac{1}{2} \text{tr}(y_m^\dagger y_n) \quad (3.69)$$

while

$$L_{mm}(x = X_n) = |X_m - X_n|^2 + \hat{a}_m^2 \approx |X_m - X_n|^2, \quad (3.70)$$

hence

$$\left[\frac{L_{mn}L_{nm}}{L_{mm}} \right] (x = X_n) \approx \frac{\hat{a}_m^2 \hat{a}_n^2 \text{tr}^2(y_m^\dagger y_n)}{4|X_m - X_n|^2} \quad (3.71)$$

and therefore

$$\mathcal{I}_{mn}^{(2)}(x \text{ near } X_n) \approx \frac{3\hat{a}_m^2 \text{tr}^2(y_m^\dagger n_j)}{\pi^2|X_m - X_n|^2} \times \frac{\hat{a}_n^4(\hat{a}_n^2 - r_n^2)}{(\hat{a}_n^2 + r_n^2)^5} = \frac{\hat{a}_m^2 \text{tr}^2(y_m^\dagger y_n)}{2|X_m - X_n|^2} \times \frac{\hat{a}_n^2 - r_n^2}{\hat{a}_n^2 + r_n^2} \times \mathcal{I}_n^{(1)}(x). \quad (3.72)$$

It is easy to check that such two-instanton interference terms do not affect the net charge of the n^{th} instanton —

$$\int_{\text{near } X_n} d^4x \mathcal{I}_{mn}^{(2)}(x) = 0 \quad (3.73)$$

— but they move the charge distribution close to the center X_n since $\mathcal{I}_{mn}^{(2)}$ is positive for $r_n < \hat{a}_n$ but negative for $r_n > \hat{a}_n$. Consequently, they decrease the effective charge radius ρ_n , which corresponds to $\Delta_n > 0$ and hence positive extra Coulomb energy. Specifically,

$$\begin{aligned} \Delta_i [\text{due to } \mathcal{I}_{mn}^2] &= \iint_{\text{instanton } \#n} d^4x d^4x' \frac{2\mathcal{I}_n^{(1)}(x) \times \mathcal{I}_{mn}^{(2)}(x')}{|x - x'|^2} \\ &= 144 \frac{\hat{a}_m^2 \text{tr}^2(y_m^\dagger y_n)}{|X_m - X_n|^2} \times \iint_0^\infty dr r^3 dr' r'^3 \frac{\hat{a}_n^4}{(\hat{a}_n^2 + r^2)^4} \times \frac{\hat{a}_n^4(\hat{a}_n^2 - r'^2)}{(\hat{a}_n^2 + r'^2)^5} \times \frac{1}{\max(r^2, r'^2)} \\ &= + \frac{\hat{a}_m^2}{5\hat{a}_n^2} \times \frac{\text{tr}^2(y_m^\dagger y_n)}{|X_n - X_j|^2}. \end{aligned} \quad (3.74)$$

Note however that this Δ_n is a correction of the instanton's charge radius or (rather $1/\rho_n^2$) *starting from the standalone radius \hat{a}_n instead of the original a_n* . Consequently, there is an additional correction

$$\Delta_n^{\text{extra}} = \frac{4/5}{\hat{a}_n^2} - \frac{4/5}{a_n^2} \approx -\frac{4/5}{a_n^4} \times (\hat{a}_n^2 - a_n^2) = -\frac{4/5}{a_n^4} \times \sum_{m \neq n} |\alpha_{mn}^\mu|^2. \quad (3.75)$$

The α_{mn}^μ are spelled out in eqs. (3.35); to the leading order

$$\alpha_{mn}^\mu \approx \alpha_{\mu mn}^{(1)} = a_m a_n \times \frac{\eta_{\mu\nu}^a (X_m - X_n)_\nu}{|X_m - X_n|^2} \times \frac{1}{2} \text{tr}(y_m^\dagger y_n (-i\tau^a)) \quad (3.76)$$

Using $\eta_{\mu\nu}^{(a)} \eta_{\mu\lambda}^{(b)} = \delta^{ab} \delta_{\nu\lambda}$, we obtain

$$|\alpha_{\mu mn}^{(1)}|^2 \approx \frac{a_m^2 a_n^2}{|X_m - X_n|^2} \times \left(\sum_a \frac{1}{4} \text{tr}^2(y_m^\dagger y_n (-i\tau^a)) = 1 - \frac{1}{4} \text{tr}^2(y_m^\dagger y_n) \right) \quad (3.77)$$

and hence

$$\Delta_n^{\text{extra}} = -\frac{1}{5} \sum_{m \neq n} \frac{a_m^2 (4 - \text{tr}^2(y_m^\dagger y_n))}{a_n^2 |X_m - X_n|^2} + O(a^2/D^4). \quad (3.78)$$

Combining the corrections (3.74) and (3.78), we obtain

$$\Delta_n^{\text{net}} = +\frac{2}{5} \sum_{m \neq n} \frac{a_m^2}{a_n^2} \times \frac{\text{tr}^2(y_m^\dagger y_n) - 2}{|X_m - X_n|^2} + O(a^2/D^4) \quad (3.79)$$

and therefore

$$\mathcal{E}_C^{\text{net}} = \frac{N_c}{4\lambda M} \left(\sum_n \frac{4/5}{a_n^2} + \sum_{m \neq n} \frac{1}{|X_m - X_n|^2} \times \left(1 + \frac{1}{5} \left(\frac{a_m^2}{a_n^2} + \frac{a_n^2}{a_m^2} \right) \times (\text{tr}^2(y_m^\dagger y_n) - 2) \right) \right) + O(a^2/D^4). \quad (3.80)$$

Finally, let us combine the non-abelian and the Coulomb energies of the system and re-organize the net energy into one-body, two-body, etc., terms:

$$\mathcal{E}^{\text{total}} = \sum_n \mathcal{E}^{\text{1 body}}(n) + \frac{1}{2} \sum_{m \neq n} \mathcal{E}^{\text{2 body}}(m, n) + \frac{1}{6} \sum_{\substack{\text{different} \\ \ell, m, n}} \mathcal{E}^{\text{3 body}}(\ell, m, n) + \dots \quad (3.81)$$

where

$$\mathcal{E}^{\text{1 body}}(n) = N_c M \left(\lambda M^2 \times (X_n^4)^2 + \frac{\lambda M^2}{2} \times a_n^2 + \frac{1}{5\lambda M^2} \times \frac{1}{a_n^2} \right), \quad (3.82)$$

$$\mathcal{E}^{\text{2 body}}(m, n) = \frac{N_c}{2\lambda M} \times \frac{1}{|X_m - X_n|_{4D}^2} \left(\begin{aligned} & \lambda^2 M^4 \times a_m^2 a_n^2 \times \text{tr}^2(y_m^\dagger y_n (-i\vec{N}_{mn} \cdot \vec{\tau})) \\ & + 1 + \frac{1}{5} \left(\frac{a_m^2}{a_n^2} + \frac{a_n^2}{a_m^2} \right) \times (\text{tr}^2(y_m^\dagger y_n) - 2) \\ & + O\left(\frac{a^2}{D^2} \sim \frac{1}{\lambda M^2 D^2}\right) \end{aligned} \right), \quad (3.83)$$

$$\mathcal{E}^{\text{3 body}}(\ell, m, n) = O\left(\frac{N_c a^2}{\lambda M^2 D^4} \sim \frac{N_c}{\lambda^2 M^3 D^4}\right), \quad \text{etc., etc.} \quad (3.84)$$

When comparing magnitudes of the non-abelian and the Coulomb terms here we have used $a \sim 1/(\sqrt{\lambda}M)$. In fact, we may be more precise by noticing that for small instantons distant from each other, the one-body potential energies for the instanton radii a_i are much larger than the two-body, etc., interactions between different instantons. Consequently, in the minimal-energy or near-minimal-energy configuration of the multi-instanton system, the instanton radii will be close to the equilibrium radius of a stand-alone instanton,

$$a_n = a_0 + O(a^3/D^2) \quad (3.85)$$

where

$$a_0 = \frac{\sqrt[4]{2/5}}{\sqrt{\lambda}M} \quad (3.86)$$

minimizes the one-body potential energy (3.82). Plugging this equilibrium radius into eq. (3.83), we obtain a simpler formula for the two-instanton interaction energy:

$$\mathcal{E}^{\text{2 body}}(m, n) = \frac{2N_c}{5\lambda M} \times \frac{1}{|X_m - X_n|_{4D}^2} \times \left[\frac{1}{2} + \text{tr}^2(y_m^\dagger y_n) + \text{tr}^2(y_m^\dagger y_n (-i\vec{N}_{mn} \cdot \vec{\tau})) \right]. \quad (3.87)$$

Note that the expression inside ‘ $[\cdots]$ ’ is always positive, so the two-body forces between the instantons are always repulsive, regardless of the instantons’ $SU(2)$ orientations. However, the orientations do affect the strength of the repulsion: two instantons with similar orientations repel each other 9 times stronger than the instantons at the same distance from each other but whose orientations differ by a 180° rotation (in $SO(3)$ terms) around a suitable axis. *This fact will be at the core of our analysis of instanton crystals in subsequent sections.*

4 Linear chains of instantons

Consider an infinite set of instantons arranged in a crystalline lattice. In light of eq. (3.87), the orientations of instantons at different lattice sites affects the system’s energy just as much as the lattice geometry. In this section we shall see how this works for a simple one-dimensional low-density lattice — i.e. an infinite chain of small instantons located along a straight line at

$$X_n^\mu = (nD, 0, 0, 0), \quad n \in \mathbb{Z} \quad (4.1)$$

where the lattice spacing D is much larger than the instanton radius a .

Before we go any further, we need a reason why the instantons would form a 1D chain rather than spread out into all 3 flat space dimensions like holographic baryons (or real-life baryons) in the compressed nuclear matter. In principle, we may sidestep this question by simply freezing the X_n^μ moduli of the instantons and allow only their orientation moduli y_n to vary, but then it would be hard to place such a system in any kind of a physical context. Alternatively, we can make 5D gauge coupling depend on the x_2 and x_3 coordinates as well as the holographic dimension x_4

$$\frac{8\pi^2}{g_5^2(x_2, x_3, x_4)} = N_c \lambda M \left(1 + M_4^2 x_4^2 + M_3^2 x_3^2 + M_2^2 x_2^2 + O(M^4 x^4) \right), \quad M_4 \equiv M, \quad (4.2)$$

which creates an effective potential for the instanton centers X_n ,

$$\mathcal{E}^{(1)}(n) = N_c \lambda M M_4^2 \times (X_n^4)^2 + N_c \lambda M M_3^2 \times (X_n^3)^2 + N_c \lambda M M_2^2 \times (X_n^2)^2 + O(N_c \lambda M^5 X^4). \quad (4.3)$$

At low instanton densities, this potential makes the instantons line up along the x^1 axis, while at higher densities, the repulsion between the instantons becomes stronger than this potential and they spread out into the other dimensions. We shall address such spreading out in the following section 5, but for the moment let us focus on the low-density 1D chains (4.1).

In a bigger holographic picture, a g_5 that depends on three coordinates $x_{2,3,4}$ indicates that the flavor branes are curved in all 3 of these directions,

$$U[\text{the energy scale}] \approx U_0 + \frac{X_4^2}{2R_4} + \frac{X_3^2}{2R_3} + \frac{X_2^2}{2R_2} + O(X^4/R^3), \quad (4.4)$$

which is quite different from the usual holographic setup where the energy-scale coordinate U depends only on the x^4 . But in this article we are not going to explore the geometric

aspects of such curvature. Instead, we shall simply use the gauge coupling (4.2) in an approximately flat space (for distances $\ll 1/M$) as a tool to put instantons into a 1D lattice. In our next article [58], we shall make g_2 depend only on x_3 and x_4 to put instantons into a 2D lattice in the (x_1, x_2) plane, and eventually go back to g_5 depending only on the x_4 to make 3D instanton lattices.

For simplicity, let us assume that the g_5 is much less sensitive to the x_2 and x_3 coordinates than the x_4 , thus $M_2^2, M_3^3 \ll M_4^2$. In this limit, the M_2^2 and M_3^3 parameters give rise to the x_2^2 and x_3^3 terms in the one-body instanton potential (4.3), but their effect on the instanton radius a or the two-body forces between the instantons may be neglected (compared to the effect of the $M_4^2 x_4^2$ term). Therefore, the net two-body forces between the instantons remain approximately as in eq. (3.87).

Moreover, for the 1D lattice geometry (4.1) of the instanton centers, we have $|X_m - X_n|^2 = D^2 \times (m-n)^2$ while $\vec{N}_{mn} = (\pm 1, 0, 0)$. Consequently, the net energy of the instanton chain as a function of the instantons' orientations y_n may be summarized as

$$\mathcal{E}^{\text{int}} = \frac{N_c}{5\lambda M D^2} \times \sum_{m \neq n} \frac{1}{(m-n)^2} \times \left[\frac{1}{2} + \text{tr}^2(y_m^\dagger y_n) + \text{tr}^2(y_m^\dagger y_n (-i\tau_1)) \right]. \quad (4.5)$$

To minimize this energy, each pair of instantons m and n wants to have $y_m^\dagger y_n$ to be a linear combination of $i\tau_2$ and $i\tau_3$; in $\text{SO}(3)$ terms, this corresponds to a relative rotation (between the 2 instantons) through a 180° angle around some axis $\perp x_1$. Alas, this cannot be achieved for all instanton pairs at once; indeed, if we minimize the energies of the (n, m) and the (n, ℓ) pairs, then the energy of the (m, ℓ) pair would be maximal instead of minimal:

$$\begin{aligned} \text{If } y_n^\dagger y_m &= i\vec{n}_1 \cdot \vec{\tau} \quad \& \quad y_n^\dagger y_\ell = i\vec{n}_2 \cdot \vec{\tau} \quad \text{for } \vec{n}_1, \vec{n}_2 \perp x_1 \\ \text{Then } y_m^\dagger y_\ell &= (\vec{n}_1 \cdot \vec{\tau})(\vec{n}_2 \cdot \vec{\tau}) = (\vec{n}_1 \cdot \vec{n}_2) + i(\vec{n}_1 \times \vec{n}_2) \cdot \vec{\tau} \quad \text{for } \vec{n}_1 \times \vec{n}_2 \parallel x_1 \\ \implies \text{tr}^2(y_m^\dagger y_\ell) + \text{tr}^2(y_m^\dagger y_\ell (-i\tau_1)) &= 4(\vec{n}_1 \cdot \vec{n}_2)^2 + 4(\vec{n}_1 \times \vec{n}_2)_1^2 = 4 \text{ (maximum)}. \end{aligned} \quad (4.6)$$

So our best bet is to first minimize the energies of the most expensive pairs — the nearest neighbors $(m, m+1)$ — and then worry about the less expensive pairs of instantons. Thus, we want

$$\forall m : \quad y_m^\dagger y_{m+1} = \cos \psi_m \times (i\tau_3) + \sin \psi_m \times (i\tau_2) \quad \text{for some angle } \psi_m, \quad (4.7)$$

and the most general solution for this set of equations (modulo a common $\text{SU}(2)$ symmetry of all the y_m) is

$$y_n = \exp(i\phi_n \tau_1) \times (i\tau_3)^n = \begin{cases} \pm [\cos \phi_n \times 1 + \sin \phi_n \times (i\tau_1)] & \text{for even } n, \\ \pm [\cos \phi_n \times (i\tau_3) + \sin \phi_n \times (i\tau_2)] & \text{for odd } n, \end{cases} \quad (4.8)$$

for some angles ϕ_n . Indeed, for any set of the angles ϕ_n we have

$$\begin{aligned} y_m^\dagger y_{m+1} &= (-i\tau_3)^m \times \exp(i(\phi_{m+1} - \phi_m)\tau_1) \times (i\tau_3)^{m+1} \\ &= \exp((-1)^m i(\phi_{m+1} - \phi_m)\tau_1) \times (i\tau_3) \\ &= \cos(\phi_{m+1} - \phi_m) \times (i\tau_3) + (-1)^m \sin(\phi_{m+1} - \phi_m) \times (i\tau_2), \end{aligned} \quad (4.9)$$

which agrees with eqs. (4.7) for $\psi_m = (-1)^m(\phi_{m+1} - \phi_m)$. Clearly, any set of ψ_m can be obtained for suitable ϕ_n , so eqs. (4.8) indeed describe all the solutions to the eqs. (4.7).

Eqs. (4.8) for various angles ϕ_n define a big family of instanton configurations. Surprisingly, all these configurations have exactly the same net energies! Indeed, for all sets of ϕ_n , all instanton pairs (m, n) with odd $m - n$ have minimal energies while pairs with even $m - n$ have maximal energies:

$$\begin{aligned}
 &\text{for odd } m - n, \quad y_m^\dagger y_n = \cos(\phi_n - \phi_m) \times (\pm i\tau_3) + \sin(\phi_n - \phi_m) \times (\pm i\tau_2) \\
 &\quad \Rightarrow \frac{1}{2} + \text{tr}^2(y_m^\dagger y_n) + \text{tr}^2(y_m^\dagger y_n(-i\tau_1)) = \\
 &\quad = \frac{1}{2} + 0 + 0 = \frac{1}{2}, \\
 &\text{for even } m - n, \quad y_m^\dagger y_n = \cos(\phi_n - \phi_m) \times (\pm 1) + \sin(\phi_n - \phi_m) \times (\pm i\tau_1) \\
 &\quad \Rightarrow \frac{1}{2} + \text{tr}^2(y_m^\dagger y_n) + \text{tr}^2(y_m^\dagger y_n(-i\tau_1)) = \\
 &\quad = \frac{1}{2} + 4\cos^2(\phi_n - \phi_m) + 4\sin^2(\phi_n - \phi_m) = \frac{9}{2}.
 \end{aligned} \tag{4.10}$$

Consequently, regardless of the angles ϕ_m , the net energy per instanton of the 1D lattice is

$$\begin{aligned}
 \mathcal{E}_{\text{per instanton}}^{\text{interaction}} &= \frac{N_c}{5\lambda MD^2} \times \sum_{m \neq n}^{\text{fixed } n} \frac{1}{(m - n)^2} \times \left[\frac{1}{2} + \text{tr}^2(y_m^\dagger y_n) + \text{tr}^2(y_m^\dagger y_n(-i\tau_1)) \right] \\
 &= \frac{N_c}{5\lambda MD^2} \times \sum_{\ell=m-n \neq 0} \frac{1}{\ell^2} \times \begin{cases} \frac{1}{2} & \text{for odd } \ell, \\ \frac{9}{2} & \text{for even } \ell, \end{cases} \\
 &= \frac{N_c}{5\lambda MD^2} \times \left(\frac{1}{2} \sum_{\text{odd } \ell} \frac{1}{\ell^2} + \frac{9}{2} \sum_{\text{even } \ell \neq 0} \frac{1}{\ell^2} \right) \\
 &= \frac{N_c}{5\lambda MD^2} \times \left(\frac{1}{2} \times \frac{\pi^2}{4} + \frac{9}{2} \times \frac{\pi^2}{12} = \frac{\pi^2}{2} \right).
 \end{aligned} \tag{4.11}$$

Now let's consider the instantons' orientations y_n from the group-theoretical point of view. Each y_n is an SU(2) matrix, but its overall sign is irrelevant, so we only care about the y_n (modulo sign), which belongs to the SU(2)/ $\mathbb{Z}_2 \cong \text{SO}(3)$. In a generic lowest-energy configuration (4.8) of the 1D chain, the instanton orientations span an $\text{SO}(2) \times \mathbb{Z}_2$ subgroup of the SO(3) corresponding to the rotational symmetries of a cylinder — rotations through arbitrary angles around the x_1 axis, and 180° rotations around axes $\perp x_1$. The y_n alternate between the two types of rotations, but apart from that they generically do not follow any regular patterns.

However, the family (4.8) also contains some regular patterns in which the orientations y_n (modulo sign) follow a repeating cycle of finite length p ; moreover, the values of y_n span a discrete subgroup of the cylindrical symmetry $\text{SO}(2) \times \mathbb{Z}_2$. Here are some examples:

- The *anti-ferromagnetic chain*, with 2 alternating instanton orientations:

$$y_{\text{even } n} = \pm 1, \quad y_{\text{odd } n} = \pm i\tau_3. \tag{4.12}$$

In this configuration — which obtains for $\phi_n \equiv 0$ — the y_n (modulo sign) span a \mathbb{Z}_2 subgroup of the $\text{SO}(2) \times \mathbb{Z}_2$.

- Period = 4 configuration spanning the *Klein group* of 180° rotations around the 3 Cartesian axes:⁶

$$y_{n \equiv 0 \pmod{4}} = \pm 1, \quad y_{n \equiv 1 \pmod{4}} = \pm \tau_3, \quad y_{n \equiv 2 \pmod{4}} = \pm \tau_1, \quad y_{n \equiv 3 \pmod{4}} = \pm \tau_2. \quad (4.13)$$

- Period = $2k = 6, 8, 10, \dots$ configurations spanning *prismatic groups* $\mathbb{Z}_k \times \mathbb{Z}_2$:

$$\begin{aligned} y_{\text{even } n} &= \cos \frac{\pi n}{2k} \times 1 + \sin \frac{\pi n}{2k} \times (i\tau_1), \\ y_{\text{odd } n} &= \cos \frac{\pi(n-1)}{2k} \times (i\tau_3) + \sin \frac{\pi(n-1)}{2k} \times (i\tau_2). \end{aligned} \quad (4.14)$$

- Period = $2k = 6, 8, 10, \dots$ configurations spanning *dihedral groups* D_{2k} , which obtain for $\phi_n = n \times (\pi/2k)$, i.e.

$$\begin{aligned} y_{\text{even } n} &= \cos \frac{\pi n}{2k} \times 1 + \sin \frac{\pi n}{2k} \times (i\tau_1), \\ y_{\text{odd } n} &= \cos \frac{\pi n}{2k} \times (i\tau_3) + \sin \frac{\pi n}{2k} \times (i\tau_2). \end{aligned} \quad (4.15)$$

There is a wider class of regular configurations — we shall call them *link-periodic* — in which the y_n themselves are not periodic, but the relative rotations $y_n^\dagger y_{n+1}$ between nearest neighbors follow a periodic pattern. In terms of the ϕ_n angles, this corresponds to periodic differences $\phi_{n+1} - \phi_n$, for example

$$\phi_n = n\varphi - \frac{1}{2}(-1)^n\theta \quad \implies \quad \phi_{n+1} - \phi_n = \begin{cases} \varphi + \theta & \text{for even } n, \\ \varphi - \theta & \text{for even } n. \end{cases} \quad (4.16)$$

For rational φ/π this pattern produces a periodic array of instantons' orientations y_n — for example, the dihedral cycle for $\varphi = \pi/2k$ and $\theta = 0$, or the prismatic cycle for $\varphi = \pi/2k$ and $\theta = -\varphi/2$. But for irrational φ 's the orientations y_n themselves do not have a finite period; instead, they wind irrationally around the cylinder group $\text{SU}(2) \times \mathbb{Z}_2$.

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All in all, the 1D lattice of instantons has a huge degenerate family (4.8) of lowest-energy configurations. However, this degeneracy obtains only in the $M_2, M_3 \ll M_4$ limit where the 5D gauge coupling (4.2) depends mostly on the holographic coordinate x_4 and only a little bit on the x_2 and x_3 . For the larger $M_2, M_3 \sim M_4$, the degeneracy is lifted and there is a unique lowest-energy configuration of the 1D lattice, namely the link-periodic array (4.16) whose φ and θ parameters depends on the M_2/M_3 ratio.

To see how this works we note that for $M_2 \sim M_3 \sim M_4$, all 3 coordinates $x_{2,3,4}$ transverse to the 1D lattice play similar roles, and the 3 terms $M_4^2 x_4^2 + M_3^2 x_3^2 + M_2^2 x_2^2$

⁶As a subgroup of $\text{SO}(3)$, the Klein group is abelian and isomorphic to $\mathbb{Z}_2 \times \mathbb{Z}_2$. But its covering group in $\text{SU}(2)$ is non-abelian group and isomorphic to the group of unit quaternions $\pm 1, \pm i, \pm j, \pm k$.

inside ‘ (\dots) ’ in eq. (4.2) give rise to similar contributions to the non-abelian energy of a multi-instanton system:

$$\mathcal{E}_{NA}^{\text{net}} = N_c \lambda M \times \int d^4x I(x) \times \sum_{\mu=2,3,4} M_\mu^2 (x_\mu)^2 \quad (4.17)$$

(where M_4^2 is the same as M^2). Consequently, the two-body interaction energy between the instantons becomes more complicated than in eq. (3.42). Instead, generalizing eq. (3.39) we obtain

$$\mathcal{E}_{NA}^{\text{net}} = N_c \lambda M \sum_{\mu=2,3,4} M_\mu^2 \left\{ \sum_n \left((X_n^\mu)^2 + \frac{1}{2} a_n^2 \right) + \sum_{n \neq m} (\alpha_{nm}^\mu)^2 \right\} \quad (4.18)$$

and hence to the leading order in expansion (3.35),

$$\mathcal{E}_{NA}^{\text{net}} \approx \sum_n \mathcal{E}_{NA}^{(1)}(n) + \frac{1}{2} \sum_{m \neq n} \mathcal{E}_{NA}^{(2)}(m, n) \quad (4.19)$$

where

$$\mathcal{E}_{NA}^{(1)}(n) = \sum_{\mu=2,3,4} N_c \lambda M M_\mu^2 (X_n^\mu)^2 + \frac{1}{2} N_c \lambda M (M_4^2 + M_3^2 + M_2^2) \times a_n^2 \quad (4.20)$$

and

$$\mathcal{E}_{NA}^{(2)}(m, n) = \frac{a_n^2 a_m^2}{2 |X_m - X_n|^2} \times \sum_{\mu=2,3,4} N_c \lambda M M_\mu^2 \times \left(\eta_{\mu\nu}^a N_{mn}^\nu \text{tr}(y_m^\dagger y_n (-i\tau^a)) \right)^2. \quad (4.21)$$

Note that the coefficient of each a_n^2 in the one-body potential (4.20) depends on the M_3^2 and the M_2^2 as well as the $M_4^2 \equiv M^2$. Consequently, when we combine this potential with the Coulomb one-body potential for the instanton radius a , the equilibrium radius ends up smaller than in eq. (3.86), namely

$$a_0^4 = \frac{2/5}{\lambda^2 M^2 (M_4^2 + M_3^2 + M_2^2)}. \quad (4.22)$$

Plugging this radius into eq. (4.21) and combining with the Coulomb two-body interaction term from eq. (3.80), we obtain the net two-body force in the background with $M_2 \sim M_3 \sim M_4 \equiv M$:

$$\mathcal{E}_{\text{net}}^{(2)}(m, n) = \frac{2N_c}{5\lambda M} \times \frac{1}{|X_m - X_n|^2} \times \left(\frac{1}{2} + \text{tr}^2(y_m^\dagger y_n) + \sum_{\mu=2,3,4} C_\mu \times \left(\eta_{\mu\nu}^a N_{mn}^\nu \text{tr}(y_m^\dagger y_n (-i\tau^a)) \right)^2 \right) \quad (4.23)$$

where

$$C_\mu \stackrel{\text{def}}{=} \frac{M_\mu^2}{M_4^2 + M_3^2 + M_2^2}, \quad C_4 + C_3 + C_2 = 1. \quad (4.24)$$

For instantons located along a straight line, this formula can be simplified a bit using $N_{mn}^\nu \equiv (\pm 1, 0, 0, 0)$ and hence

$$\eta_{\mu\nu}^a N_{mn}^\nu \operatorname{tr}(y_m^\dagger y_n (-i\tau^a)) = \begin{cases} \pm \operatorname{tr}(y_m^\dagger y_n (-i\tau^1)) & \text{for } \mu = 4, \\ \pm \operatorname{tr}(y_m^\dagger y_n (-i\tau^2)) & \text{for } \mu = 3, \\ \pm \operatorname{tr}(y_m^\dagger y_n (-i\tau^3)) & \text{for } \mu = 2. \end{cases} \quad (4.25)$$

Consequently, the net interaction energy of the 1D lattice is

$$\mathcal{E}_{\text{net}}^{\text{int}} \equiv \frac{1}{2} \sum_{n \neq m} \mathcal{E}^{(2)}(n, m) = \frac{N_c}{5\lambda M D^2} \times \sum_{m \neq n} \frac{Q(m, n)}{(m - n)^2} \quad (4.26)$$

where

$$Q(m, n) \stackrel{\text{def}}{=} \frac{1}{2} + \operatorname{tr}^2(y_m^\dagger y_n) + C_4 \operatorname{tr}^2(y_m^\dagger y_n (-i\tau^1)) \\ + C_3 \operatorname{tr}^2(y_m^\dagger y_n (-i\tau^2)) + C_2 \operatorname{tr}^2(y_m^\dagger y_n (-i\tau^3)). \quad (4.27)$$

Without loss of generality we assume $M_4 \geq M_3 \geq M_2$ and hence $C_4 \geq C_3 \geq C_2$ — otherwise we simply re-label the coordinates $x_{2,3,4}$. Thus, the lowest-energy relative orientation of an instanton pair (m, n) is $y_m^\dagger y_n = \pm i\tau^3$. If this orientation is unachievable, the next best bet is a linear combination of $i\tau^3$ and $i\tau^2$. If that is also unachievable, the third-best choice would be a linear combination of all three $i\tau^{1,2,3}$; in $\text{SO}(3)$ terms, this corresponds to the 180° rotation around a generic axis. Finally, rotations through other angles would be the least attractive option which the instantons $\#m$ and $\#n$ would rather avoid — unless they are forced to it by interactions with the other instantons between m and n .

In this setting, it is not intuitively obvious how to balance the energy cost of nearest-neighbor interactions versus next-to-nearest neighbors and more distant instanton pairs. Instead of intuition, we have performed a computer experiment using a lattice of 200 $\text{SU}(2)$ matrices y_n . In each run, we start with the y_n being independent random elements of the $\text{SU}(2)$ group, and then let them evolve towards a minimum of the energy function (4.26) via the relaxation method. That is, we let the y_n evolve with time according to

$$\frac{dy_n(t)}{dt} = -K \times \frac{\delta \mathcal{E}_{\text{net}}^{\text{int}}}{\delta y_n} \quad (4.28)$$

where K is a constant mobility factor and the derivative with respect to an $\text{SU}(2)$ matrix y_n is defined as

$$\frac{\delta \mathcal{E}}{\delta y_n} \stackrel{\text{def}}{=} y_n \left((-i\vec{\tau}) \cdot \nabla_s \mathcal{E}(y_n \rightarrow y_n(1 + i\vec{s} \cdot \vec{\tau})) \Big|_{\vec{s}=0} \right). \quad (4.29)$$

Each run ends when all the y_n seem to converge to an equilibrium configuration and their derivatives (4.28) become very small. We made many runs for different C_3 and C_4 parameters, and here is what we have found:

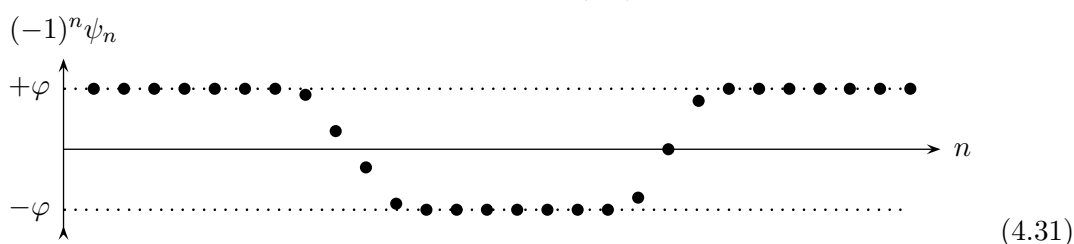
- In all equilibrium configurations for backgrounds with $C_3 < C_4$, all twists $y_n^\dagger y_{n+1}$ between nearest neighbors have form

$$y_n^\dagger y_{n+1} = \cos \psi_n \times i\tau^3 + \sin \psi_n \times i\tau^2 + (\text{tiny}) \times i\tau^1 + (\text{tiny}) \times 1, \quad (4.30)$$

where the tiny coefficients of 1 and $i\tau^1$ are artefacts of imperfect convergence to equilibrium (they get smaller when we allow more time for convergence).

- The backgrounds with $M_3 = M_4 \implies C_3 = C_4$ are trickier due to symmetry between $i\tau^1$ and $i\tau^2$ components of the $SU(2)$ matrices. (Or all three $i\tau^{1,2,3}$ matrices for $M_2 = M_3 = M_4 \implies C_2 = C_3 = C_4$.) In equilibrium configurations for such symmetric backgrounds, the nearest-neighbor twists $y_n^\dagger y_{n+1}$ generally do have $i\tau^1$ components. However, all such components can be eliminated by a suitable global symmetry of the system, $y_n \rightarrow y_n \times G$, same $G \in SU(2)$ for all y_n , and then all the nearest-neighbor twists $y_n^\dagger y_{n+1}$ take the form (4.30).
- For most nearest-neighbor pairs, the angles ψ_n in eqs. (4.30) take values $\pm\varphi \pmod{\pi}$ where φ depends on the $C_{2,3,4}$ parameters. Moreover, the \pm sign here tends to alternate between odd and even n . However, some pairs do not follow these rules.

In general, our runs end up with patterns of $(-1)^n \psi_n$ that look like this:



Physically, this corresponds to the 1D lattice having two degenerate ground states, one with $\psi_n = +(-1)^n \varphi$ and the other with $\psi_n = -(-1)^n \varphi$. In our numerical runs, some parts of the lattice converge to one of these states while other parts converge to the other state; altogether, we end up with several domains separated by walls. In a perfect simulation, the walls would eventually move towards each other and annihilate, so the whole lattice would end up in the same ground state. But this process is so slow that our numeric simulation stops before it barely begins, thus we always end up with multiple domains instead of a single ground state for the whole lattice.

- ★ The bottom line of our numerical simulations is that the ground state of the 1D instanton lattice always has a link-periodic instanton orientations (4.16) (up to a global symmetry, if any) for a periodicity angle φ that depends on ratios of parameters $M_2 : M_3 : M_4$. Curiously, it actually depends only on the $M_2 : M_3$ ratio regardless of the M_4 (as long as $M_4 \geq M_3 \geq M_2$).

Once we know the ground state of the instanton orientations in a 1D lattice is link-periodic, we may calculate the periodicity angle φ analytically. Indeed, applying eqs. (4.10)

to the link-periodic array (4.16) of instanton orientations, we have:

$$\begin{aligned}
 \text{For even } n-m, \quad y_m^\dagger y_n &= \cos((n-m)\varphi) \times i^{n-m} + \sin((n-m)\varphi) \times i^{n+m+1}\tau^1 \\
 \implies Q(m, n) &= \frac{1}{2} + 4\cos^2((n-m)\varphi) + 4C_4 \sin^2((n-m)\varphi) \\
 &= \frac{5}{2} + 2C_4 + 2(1 - C_4 = C_3 + C_2) \times \cos(2(n-m)\varphi),
 \end{aligned} \tag{4.32}$$

$$\begin{aligned}
 \text{For odd } n-m, \quad y_m^\dagger y_n &= \cos((n-m)\varphi \pm \theta) \times i^{n-m}\tau^3 + \sin((n-m)\varphi \pm \theta) \times i^{n+m}\tau^2 \\
 \implies Q(m, n) &= \frac{1}{2} + 4C_2 \cos^2((n-m)\varphi \pm \theta) + 4C_3 \sin^2((n-m)\varphi \pm \theta) \\
 &= \frac{1}{2} + 2(C_3 + C_2) - 2(C_3 - C_2) \times \cos(2(n-m)\varphi \pm 2\theta),
 \end{aligned} \tag{4.33}$$

where the \pm sign is $(-1)^n$. To calculate average interaction energy per instanton in the lattice, we should average $Q(n, m)$ between odd and even n for fixed $\ell = n - m$; indeed

$$\mathcal{E}_{\text{per instanton}}^{\text{interaction}} = \frac{N_c}{5\lambda MD^2} \times \left\langle \sum_{m \neq n}^{\text{fixed } n} \frac{Q(m, n)}{(n-m)^2} \right\rangle_{\text{over } n}^{\text{average}} = \frac{N_c}{5\lambda MD^2} \times \sum_{\ell \neq 0} \frac{\bar{Q}(\ell)}{\ell^2}. \tag{4.34}$$

Thus,

$$\begin{aligned}
 \text{for even } \ell, \quad \bar{Q}(\ell) &= \frac{5}{2} + 2C_4 + 2(C_3 + C_2) \cos(2\ell\varphi), \\
 \text{for odd } \ell, \quad \bar{Q}(\ell) &= \frac{1}{2} + 2(C_3 + C_2) - 2(C_3 - C_2) \times \cos(2\theta) \times \cos(2\ell\varphi),
 \end{aligned} \tag{4.35}$$

and therefore

$$\begin{aligned}
 \mathcal{E}_{\text{per instanton}}^{\text{interaction}} &= \frac{N_c}{5\lambda MD^2} \times \left(\begin{aligned} &\left(\frac{5}{2} + 2C_4 \right) \times \left(\sum_{\text{even } \ell \neq 0} \frac{1}{\ell^2} = \frac{\pi^2}{12} \right) \\ &+ 2(C_3 + C_2) \times \left(\sum_{\text{even } \ell \neq 0} \frac{\cos(2\ell\varphi)}{\ell^2} = \frac{\pi^2}{12} - \pi|\varphi| + 2\varphi^2 \right) \\ &+ \left(\frac{1}{2} + 2C_3 + 2C_2 \right) \times \left(\sum_{\text{odd } \ell} \frac{1}{\ell^2} = \frac{\pi^2}{4} \right) \\ &- 2(C_3 - C_2) \cos(2\theta) \times \left(\sum_{\text{odd } \ell} \frac{\cos(2\ell\varphi)}{\ell^2} = \frac{\pi^2}{4} - \pi|\varphi| \right) \\ &\ll \text{where the sums are evaluated assuming } |\varphi| < \frac{\pi}{2} \gg \end{aligned} \right) \\
 &= \frac{N_c}{5\lambda MD^2} \times \left(\begin{aligned} &\frac{\pi^2}{2} \times \left(1 + C_3 + C_2 - (C_3 - C_2) \cos(2\theta) \right) \\ &- 2\pi|\varphi| \times \left(C_3 + C_2 - (C_3 - C_2) \cos(2\theta) \right) \\ &+ 4\varphi^2 \times \left(C_3 + C_2 \right) \end{aligned} \right). \tag{4.36}
 \end{aligned}$$

Minimizing this expression WRT the φ and θ produces 4 degenerate minima, namely

$$\begin{aligned}
 (1) \quad & \varphi = +\frac{\pi}{2} \times \frac{C_2}{C_2 + C_3}, \quad \theta = 0, \\
 (2) \quad & \varphi = -\frac{\pi}{2} \times \frac{C_2}{C_2 + C_3}, \quad \theta = 0, \\
 (3) \quad & \varphi = -\frac{\pi}{2} \times \frac{C_3}{C_2 + C_3}, \quad \theta = \frac{\pi}{2}, \\
 (4) \quad & \varphi = +\frac{\pi}{2} \times \frac{C_3}{C_2 + C_3}, \quad \theta = \frac{\pi}{2},
 \end{aligned} \tag{4.37}$$

however the last two minima are physically equivalent to the first two. Thus, in agreement with our computer ‘experiments’, the 1D instanton lattice has 2 degenerate ground states related by the $\varphi \rightarrow -\varphi$ symmetry. The value of $|\varphi|$ also agrees with our ‘experimental data’:

$$\theta = 0, \quad \varphi = \pm \frac{\pi}{2} \times \frac{M_2^2}{M_2^2 + M_3^2} \quad \text{regardless of the } M_4^2 \tag{4.38}$$

(as long as $M_4 \geq M_3 \geq M_2$). In particular, for rational ratios of M_2^2/M_3^2 , the angle φ is rational (in units of π) and the instanton orientations y_n become periodic. Specifically, for

$$\frac{M_2^2}{M_3^2} = \frac{p}{q}, \quad p < q, \quad \gcd(p, q) = 1, \tag{4.39}$$

the y_n (modulo signs) repeat with period $2(q + p)$ while spanning the $D_{2(q+p)}$ dihedral subgroup of the $SO(3) \cong SU(2)/\mathbb{Z}_2$, for example see eq. (4.15) for $p = 1$ and $q = k - 1$.

Two particularly interesting $M_2^2 : M_3^2$ ratios need special handling, $M_2^2 = M_3^2$ and $M_2^2 \ll M_3^2$. For $M_2 = M_3$ the background has a rotational symmetry in the $x^{2,3}$ plane; for the 1D instanton lattice, this translates into the $U(1)$ symmetry between the $i\tau^2$ and $i\tau^3$ directions in the $SU(2)$. Consequently, instead of two discrete ground states (4.38) we have a continuous family:

$$\varphi = \frac{\pi}{2}, \quad \theta = \text{anything}. \tag{4.40}$$

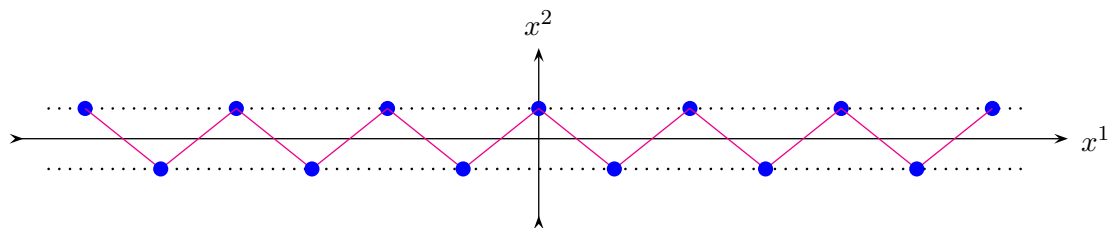
For all configurations in this family, the instanton orientations y_n (modulo signs) repeat with period 4 while spanning the Klein groups $\mathbb{Z}_2 \times \mathbb{Z}_2$ of 180° rotations around 3 mutually \perp axes. For $\theta = -\frac{\pi}{4}$ the y_n are spelled out in eq. (4.13); for other values of θ we have similar cycles related by the $U(1) \subset SU(2)$ symmetry.

Finally, for the very asymmetric background with $M_2^2 \ll M_3^2$, the two ground states (4.38) become indistinguishable as $\varphi \rightarrow \pm 0$. In this limit, the instanton orientations form the anti-ferromagnetic order (4.12).

5 Instanton zigzags

In this section we consider the most likely first step in the transition between 1D and 2D instanton lattices — a lattice with one infinite dimension x^1 while the other dimension x^2 has just two layers. Since the instantons repel each other, the two layers should be

staggered in the x^1 direction, so the whole lattice looks like a single zigzag-shaped chain:



$$(5.1)$$

In such a zigzag, the instanton $\#n$ has its center at

$$X_n^1 = n \times D, \quad X_n^2 = (-1)^n \times \epsilon, \quad X_n^3 = X_n^4 = 0, \quad (5.2)$$

for some parameters D and ϵ ; we shall refer to D as the *lattice spacing* and to ϵ as the *zigzag amplitude*.

To make the instantons form a zigzag lattice, we use the same background

$$\frac{8\pi^2}{g_5^2(x_2, x_3, x_4)} = N_c \lambda M \left(1 + M^2 x_4^2 + M_3^2 x_3^2 + M_2^2 x_2^2 + O(M^4 x^4) \right), \quad (4.2)$$

we have used in the last section to make a straight chain but increase the instanton density $\rho = 1/D$ until the repulsion between the instantons pushes them away from their neighbors in some transverse direction. For $M_2 < M_3 \leq M_4$, the initial breakout from the straight line is going to be in the $\pm x^2$ directions, with opposite signs for the nearest neighbors (to increase the distance between them), hence the zigzag geometry (5.2). For higher densities — and hence stronger repulsion between the instantons — we will get more complicated lattices: multiple layers in the x^2 direction, and eventually breaking out into the x^3 and x^4 dimensions. However, the first geometry right after the straight 1D lattice ought to be the zigzag.

Clearly, the above heuristic argument is not a proof. The proof — or disproof — will come from a computer simulation of multi-instanton systems where the instantons are allowed to move independently in all directions as well as to change their orientations. We shall describe such a simulation — and its results — in our next paper [58] in this project. Meanwhile, in the present paper we shall simply assume that the instanton centers form a zigzag-shaped chain (5.1) and focus on the instantons' orientations.

Since our formulae for the instanton interactions presume distances $|X_m - X_n|$ much larger than the instanton radius a , we need the transition from a straight 1D chain to a zigzag to happen at a critical lattice spacing $D_c \gg a$, which calls for

$$M_2 \ll M_4 \equiv M. \quad (5.3)$$

Indeed, consider an instanton that have moved a little bit away from the x^1 axis in the x^2 direction. The restoring force on this instanton due to x^2 dependence of the gauge coupling is

$$F_{\text{rest}} = -X_2 \times 2N_c \lambda M M_2^2 \quad (5.4)$$

while the repulsion from the neighboring instantons pushes it further away with the net force

$$F_{\text{rep}} = -\frac{\partial}{\partial X_2} O\left(\frac{N_c}{\lambda M} \times \frac{1}{O(D^2) + X_2^2}\right) = +\frac{X_2}{D^4} \times O\left(\frac{N_c}{\lambda M}\right). \quad (5.5)$$

At the critical lattice spacing, the two forces balance each other, hence

$$D_c^4 = \frac{O(1)}{\lambda^2 M^2 M_2^2}. \quad (5.6)$$

At the same time,

$$a^4 = \frac{2/5}{\lambda^2 M^2 (M_4^2 + M_3^2 + M_2^2)} = \frac{O(1)}{\lambda^2 M^2 M_4^2}, \quad (5.7)$$

so to assure $D_c \gg a$ we need $M_2 \ll M_4 \equiv M$.

Note however that while we need $M_2 \ll M_4$, there are no restrictions on the M_3 parameter except $M_2 \leq M_3 \leq M_4$. Consequently, we may consider zigzags in a variety of backgrounds with different M_3/M_4 ratios — ranging from $M_2 = M_3 \ll M_4$ to $M_2 \ll M_3 = M_4$ — and hence different ways in which the forces between two instantons depend on their relative orientations. For small $M_3 \ll M_4$, the two-instanton interactions are spelled out in eq. (3.87) while for finite M_3/M_4 ratios we need to use a more complicated formula (4.23) (with $C_2 \approx 0$ but finite C_3 and C_4). For a zigzag where all the instantons lie in the $x^{1,2}$ plane and hence all the N_{mn}^μ have form $N_{mn}^\mu = (*, *, 0, 0)$, we may simplify this formula using the identity

$$\sum_{\mu=2,3} \left(\eta_{\mu\nu}^\alpha N_{mn}^\nu \text{tr}(y_m^\dagger y_n (-i\tau^a)) \right)^2 = \sum_{a=1,2} \text{tr}^2(y_m^\dagger y_n (-i\tau^a)) \quad (\text{for } N_{mn}^\nu = (*, *, 0, 0)). \quad (5.8)$$

Consequently, eq. (4.23) can be rearranged as

$$\mathcal{E}_{\text{net}}^{\text{int}}[\text{zigzag}] = \frac{N_c}{5\lambda M} \sum_{m \neq n} \frac{Q_z(m, n)}{|X_m - X_n|^2} \quad (5.9a)$$

$$\begin{aligned} \text{where } Q_z(m, n) = & \frac{1}{2} + \text{tr}^2(y_m^\dagger y_n) + C_3 \sum_{a=1,2} \text{tr}^2(y_m^\dagger y_n (-i\tau^a)) \\ & + (1 - 2C_3) \times \text{tr}^2(y_m^\dagger y_n (-i\vec{\tau} \cdot \vec{N}_{mn})). \end{aligned} \quad (5.9b)$$

Note that in general, the force between two instantons is not central — it depend not only on the distance between the instantons and their relative orientation $y_m^\dagger y_n$ of their isospins, but also on the direction \vec{N}_{mn} of their separation $\vec{X}_n - \vec{X}_m$ in space. However, in the background with $M_3 = M_4$ — which is the background we have used in our previous paper [1] (albeit in different notations) — $C_3 = \frac{1}{2}$, the last term in eq. (5.9b) goes away, and the force becomes central (but isospin-dependent).

The forces between instantons will be important in our next article [58] where we shall allow them to move in the $x^{1,2}$ plane seeking the lowest-energy lattice. For the moment, we simply assume that the instantons somehow form a zigzag of some particular lattice spacing D and amplitude ϵ and focus on their orientation moduli y_n . Our immediate goal

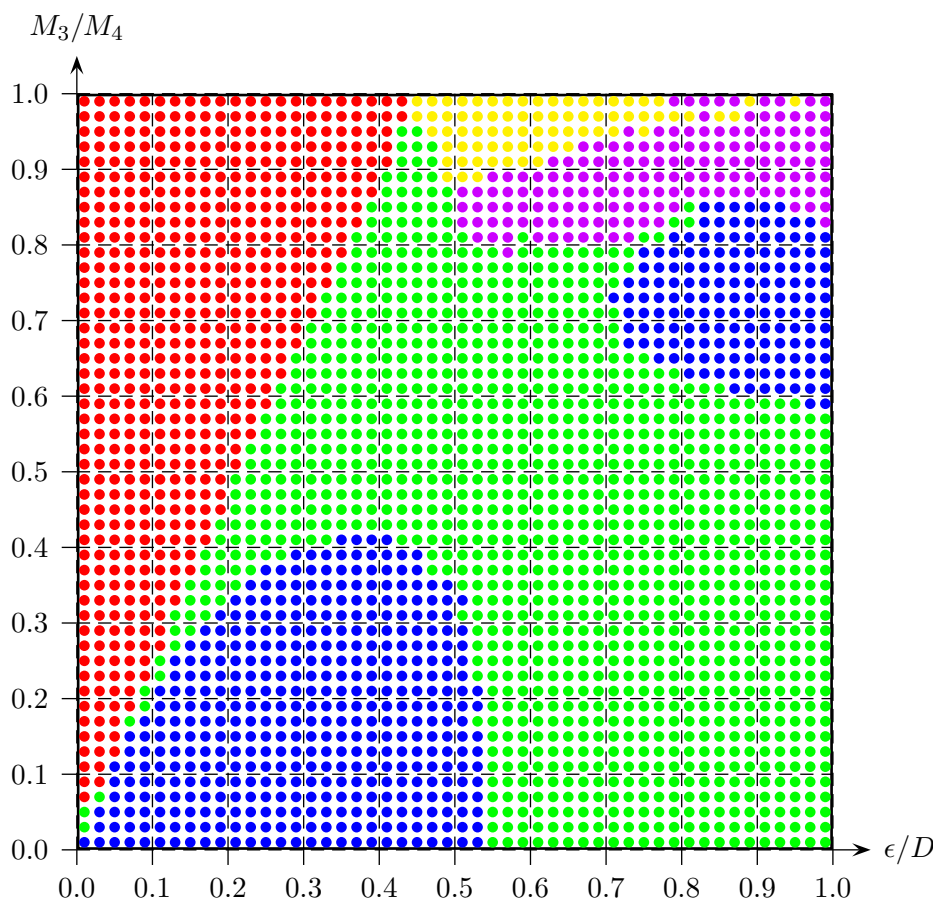


Figure 4. Rough distribution of instanton orientation patterns in the zigzag parameter space. This diagram is obtained from a numerical relaxation method.

is to find the lowest-energy configuration of the orientations y_n as a function of the ϵ/D and M_3/M_4 ratios.

A priori, we do not know if the orientations — or rather the relative orientations $y_n^\dagger y_{n+1}$ of the neighbor instantons — follow any particular pattern or patterns. To find such pattern(s) we used numerical simulations: starting with completely random y_n , we let them evolve towards a minimum of the energy function (5.9) according to eqs. (4.28); when the evolution seems to stop, we look at the relative orientations of the nearest neighbors $y_n^\dagger y_{n+1}$ to see if they form any recognizable pattern. Repeating this procedure for different combination of the ϵ/D and M_3/M_4 parameters, we saw five distinct patterns of orientations (or rather four distinct patterns and one confused mess). Figure 4 diagrams the distribution of these patterns in the parameter space.

Here is the color key to this figure:

- The red dots on figure 4 denote the anti-ferromagnetic pattern (AF) of instanton orientations in which the nearest neighbors always differ by a 180° rotation around

the third axis,

$$y_n = \begin{cases} \pm 1 & \text{for even } n, \\ \pm i\tau_3 & \text{for odd } n, \end{cases}, \quad \text{same } y_n^\dagger y_{n+1} = i\tau_3 \text{ for all } n. \quad (5.10)$$

- The yellow dots denote another abelian pattern (AB) in which all nearest neighbors differ by the same $U(1) \subset SU(2)$ rotation, but now the rotation angle is $< 180^\circ$,

$$\text{same } y_n^\dagger y_{n+1} = \exp\left(\frac{i}{2}\phi\tau_3\right) \text{ for all } n, \quad 0 < \phi < \pi. \quad (5.11)$$

- The blue dots denote a non-abelian link-periodic pattern (NA1) in which the relative rotation between nearest neighbors is always through a 180° angle, but the direction of rotation alternates between two different axes in the (12) plane, one axis for the odd-numbered instantons and the other for the even-numbered. In $SU(2)$ terms,

$$\begin{aligned} y_{2k}^\dagger y_{2k+1} &= \exp\left(\frac{i\pi}{2} \vec{n}_e \cdot \vec{\tau}\right) = i\vec{n}_e \cdot \vec{\tau} = +iA\tau_1 + iB\tau_2, \\ y_{2k+1}^\dagger y_{2k+2} &= \exp\left(\frac{i\pi}{2} \vec{n}_o \cdot \vec{\tau}\right) = i\vec{n}_o \cdot \vec{\tau} = +iA\tau_i - iB\tau_2, \end{aligned} \quad (5.12)$$

for some $A, B \neq 0$ ($A^2 + B^2 = 1$).

- The green dots denote another non-abelian link-periodic pattern (NA2). Again, the relative rotation between nearest neighbors is always through a 180° angle, but the direction of rotation alternates between two different axes. However, this time the two axes no longer lie within the (12) plane, thus

$$\begin{aligned} y_{2k}^\dagger y_{2k+1} &= iA\tau_1 + iB\tau_2 + iC\tau_3, \\ y_{2k+1}^\dagger y_{2k+2} &= iA\tau_1 - iB\tau_2 - iC\tau_3, \end{aligned} \quad (5.13)$$

where A, B, C all $\neq 0$ ($A^2 + B^2 + C^2 = 1$).

- Finally, the purple dots denote a confused non-abelian pattern (NAX) in which the nearest neighbors seem to differ by rotations through angles $\phi < 180^\circ$ around 2 alternating axes. In $SU(2)$ terms, we see something like

$$\begin{aligned} y_{2k}^\dagger y_{2k+1} &= +iA\tau_1 + iB\tau_2 + iC\tau_3 + D, \\ y_{2k+1}^\dagger y_{2k+2} &= -iA\tau_1 + iB\tau_2 + iC\tau_3 + D, \end{aligned} \quad (5.14)$$

where A, B, C, D all $\neq 0$ ($A^2 + B^2 + C^2 + D^2 = 1$),

but this picture is rather noisy: the coefficients A, B, C, D fluctuate from n to n and from run to run much stronger than their analogues in the other 4 patterns (5.10)–(5.13), and even the relative signs between A, B, C, D flip every few lattice sites — the long-range correlations are quite poor.

Physically, the disorder in the purple part of the phase diagram 4 indicates a nearby first-order phase transition between incompatible orientation patterns. Near such a transition, our relaxation method fails to converge to a homogeneous phase that happens to have lowest energy for the $(\epsilon/D, M_3/M_4)$ parameters in question. Instead, it leads to a mishmash of short domains of two nearly-degenerate phases; in fact, the domains are only a few lattice sites long, so they are significantly distorted by the boundary effects. The resulting mess is not even a local minimum of the zigzag's energy function, but once our simulation reaches this point, further evolution of the instanton orientations y_n becomes very slow. Given very long time (and a better numeric algorithm) we would eventually see the domains of each phase growing longer and eventually merging into a single phase — but in our calculations we have simply run out of time and patience before this can happen.

The bottom line of our numeric analysis is that the instanton zigzag has 4 or 5 distinct orientation phases — we can not tell if **NAX** is a real phase or an artefact of mixed domains of other phases that we could not resolve. To answer that question — and also to get a better map of phase boundaries — we need to go back to analytic work.

Our starting point is the *qualitative* result of our numerical analysis: all 5 patterns (5.10)–(5.14) we have seen — or might have seen — can be fit into a single ansatz:

$$y_n = \exp\left(in \frac{\phi}{2} \tau_2\right) \times \exp\left(i \left(\frac{\alpha}{2} + (-1)^n \frac{\beta}{2}\right) \tau_1\right), \quad (5.15)$$

$$\begin{aligned} y_n^\dagger y_{n+1} = & \cos \frac{\phi}{2} \cos \beta \times 1 - (-1)^n \cos \frac{\phi}{2} \sin \beta \times i \tau_1 \\ & + \sin \frac{\phi}{2} \cos \alpha \times i \tau_2 + \sin \frac{\phi}{2} \sin \alpha \times i \tau_3 \end{aligned} \quad (5.16)$$

for some angles ϕ, α, β . Indeed, the **AF** phase obtains for $\phi = \pi, \alpha = \frac{\pi}{2}$; the **AB** phase obtains for $0 < \phi < \pi, \alpha = \frac{\pi}{2}, \beta = 0$; the **NA1** phase obtains for $0 < \phi < \pi, \alpha = 0, \beta = \frac{\pi}{2}$; the **NA2** phase obtains for $0 < \phi < \pi, 0 < \alpha < \frac{\pi}{2}, \beta = \frac{\pi}{2}$; and the **NAX** phase — if it exists at all — obtains for generic values of all three angles, $0 < \phi < \pi, 0 < \alpha < \frac{\pi}{2}, 0 < \beta < \frac{\pi}{2}$.

Given the ansatz (5.15), we may calculate the zigzag's net energy as an analytic function of the angles ϕ, α, β and parameters $(\epsilon/D, M_3/M_4)$. The calculation is presented in appendix A, and the result is

$$\mathcal{E}_{\text{per instanton}}^{\text{zigzag}} = \frac{\pi^2 N_c}{20 \lambda M D^2} \times \mathcal{F}(\phi, \alpha, \beta; C_3, \epsilon/D) + N_c \lambda M M_2^2 \times \epsilon^2, \quad (5.17)$$

$$\begin{aligned} \text{for } \mathcal{F} = & \frac{3}{2} + \left(-1 + C_3(\cos^2 \alpha \cos^2 \beta + \sin^2 \alpha \sin^2 \beta)\right) \times \Sigma_0(\phi) \\ & + \left(\frac{5}{2} + \cos^2 \alpha - \sin^2 \beta\right) \times \Sigma_1(\epsilon/D) + \left(2 - \cos^2 \alpha - \sin^2 \beta\right) \times \Sigma_2(\phi, \epsilon/D) \\ & + (1 - 2C_3)(\sin^2 \beta - \cos^2 \alpha) \times \Sigma_3(\epsilon/D) \\ & + (1 - 2C_3)(\sin^2 \beta + \cos^2 \alpha) \times \Sigma_4(\phi, \epsilon/D) \\ & - 2(1 - 2C_3) \cos \alpha \sin \beta \times \Sigma_5(\phi, \epsilon/D), \end{aligned} \quad (5.18)$$

where

$$\Sigma_0 = \frac{4\phi(\pi - \phi)}{\pi^2}, \quad (5.19a)$$

$$\Sigma_1 = \frac{\tanh(\pi\epsilon/D)}{(\pi\epsilon/D)}, \quad (5.19b)$$

$$\Sigma_2 = \frac{\sinh((\pi - 2\phi)\epsilon/D)}{(\pi\epsilon/D) \cosh(\pi\epsilon/D)}, \quad (5.19c)$$

$$\Sigma_3 = \frac{1}{\cosh^2(\pi\epsilon/D)}, \quad (5.19d)$$

$$\Sigma_4 = \frac{\cosh(2\phi\epsilon/D)}{\cosh^2(\pi\epsilon/D)} - \frac{2\phi}{\pi} \times \frac{\cosh((\pi - 2\phi)\epsilon/D)}{\cosh(\pi\epsilon/D)}, \quad (5.19e)$$

$$\Sigma_5 = \frac{\sinh(2\phi\epsilon/D)}{\cosh^2(\pi\epsilon/D)} + \frac{2\phi}{\pi} \times \frac{\sinh((\pi - 2\phi)\epsilon/D)}{\cosh(\pi\epsilon/D)}, \quad (5.19f)$$

In appendix A we also minimize \mathcal{F} — and hence the net energy per instanton — with respect to the orientation parameters α , β , and ϕ for fixed C_3 and ϵ/D . In particular, we show that the minimum never lies at generic values of the angles ϕ, α, β — which would correspond to the **NAX** phase. This means that *the NAX phase does not really exist* — its appearance in the purple region on the figure 4 is an artefact of poor convergence near a first-order phase transition (see below).

On the other hand, for appropriate values of the ϵ/D and M_3/M_4 parameters, the zigzag energy (5.17) has global minima corresponding to the other 4 phases (5.10)–(5.13) of instanton orientations. Figure 5 on the next page shows their distributions in the parameter space. Curiously, one of the non-abelian phases — the **NA1** — appears in two different regions of the parameter space — one at low M_3/M_4 and small zigzag amplitudes, and the other at higher M_3/M_4 and higher amplitudes — while the intermediate region is occupied by the different non-abelian phase **NA2**. Both phases transitions between the two non-abelian phases are second-order. Likewise, the transition between the non-abelian **NA2** and the abelian anti-ferromagnetic phase **AF** is also second-order.

On the other hand, the transitions between the other abelian **AB** phase and the non-abelian phases **NA1** and **NA2** are first-order. In fact, these are precisely the first-order transitions that confused our numeric analysis in the purple region of the figure 4 into appearance of the fifth orientation phase **NAX**.

Finally, the triple point at $(M_3/M_4) = 1$, $(\epsilon/D) \approx 0.38$ between the **AF**, **NA2**, and **AB** phases is *critical*. At that point, the transitions between all 3 phases are second-order. Indeed, in our previous paper [1] — where the analysis was limited to zigzags with $M_3 = M_4$ (in present notations) — we saw a second-order transition between the two abelian phases **AF** and **AB**.

★ ★ ★

Thus far, we have focused on the instanton orientations y_n for a given zigzag geometry. That is, we have not only assumed that the instanton centers form a zigzag, but we have also treated the zigzag amplitude ϵ as an independent input parameter, just like

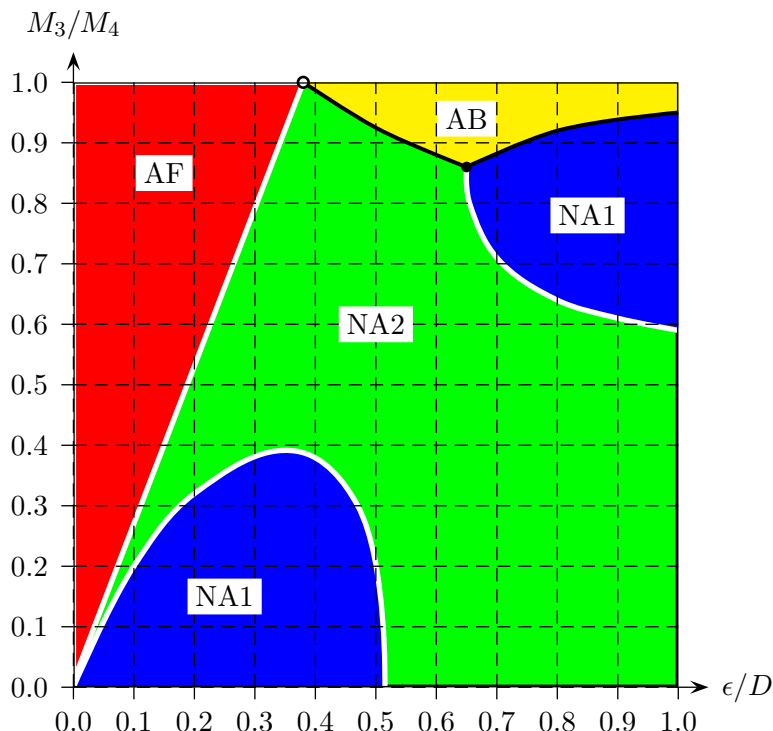


Figure 5. Phase diagram of the instanton orientation patterns in the zigzag parameter space. This diagram obtains from the mostly-analytical calculation in appendix A. The lines separating different phases indicate the order of the phase transition: a white line for a second-order transition and a black line for a first-order transition.

the lattice spacing D or the M_2, M_3, M_4 parameters of the 5D gauge theory. Indeed, figure 5 shows the transitions between different orientation phases as function of independent, freely-adjustable parameters ϵ/D and M_3/M_4 . But physically, the zigzag amplitude ϵ is a dynamical modulus whose value follows from minimizing the net energy of the multi-instanton system.

In fact, all instanton center coordinates X_n^μ are dynamical moduli, which raises two problems. First, in some situations the lowest-energy configuration of the instanton lattice may be more complicated than a zigzag or a straight chain; we shall address this issue in our next paper [58]. Second, for some lattice spacings D , a uniform lattice of any kind — a straight chain, a zigzag, or anything else — may be unstable against breaking into domains of different phases with different densities.

By way of analogy, consider a fluid governed by an equation of state such as Van-der-Waals. Formally, this equation allows a uniform fluid to have any density (up to some maximum). But in reality, at sub-critical temperatures one may have a low-density gas or a high-density liquid, but there are no uniform fluids with intermediate densities. If we constrain the overall volume V of some amount of fluid such that its *average density* would fall into the intermediate range, we would not get a uniform fluid; instead, part of the volume would be filled by the higher-density liquid while the other part by the lower-density gas. In the same way, if we fix the overall length L of the x^1 axis occupied by

some large number N of instantons, their lowest-energy configuration is not necessarily a uniform lattice of some kind. Instead, for some *average densities* $\rho = N/L$ we would have L split into domains of two different lattices of different densities, one denser than N/L and the other less dense.

To keep the fluid uniform, one should control its pressure P rather than the volume V ; consequently, the preferred phase follows from minimizing the free enthalpy $G = E - ST + PV$ rather than the free energy $F = E - ST$. Likewise, for the 1D lattice of instantons, we should control the net compression force \mathbf{F} along the x^1 axis rather than the net length L or the lattice spacing D . Also, we should minimize the free enthalpy of the lattice $G = E - ST + L\mathbf{F}$, but since we work at zero temperature all we need is the ordinary enthalpy $H = E + L\mathbf{F}$. Equivalently, we may minimize the *non-relativistic chemical potential*

$$\hat{\mu} = \mu_{\text{rel}} - M_{\text{baryon}} = \frac{G_{\text{non-rel}}}{N} \xrightarrow{T=0} \frac{E + L\mathbf{F}}{N} = \mathcal{E} + \frac{\mathbf{F}}{\rho}. \quad (5.20)$$

(We focus on the non-relativistic chemical potential $\hat{\mu} = \mu - M_{\text{baryon}}$ because the relevant scale of $\hat{\mu}$ would be *much* smaller than the baryon mass. Indeed, according to eq. (5.29) below, $\hat{\mu} \sim N_c M_2 \ll N_c M \ll N_c M \lambda \sim M_b$.)

Thus, through the remainder of this section, we are going to impose a compression force \mathbf{F} on the multi-instanton system, *assume* that the instantons form a uniform zigzag of some lattice spacing D and amplitude ϵ — or a uniform straight chain for $\epsilon = 0$ — and vary D , ϵ and the orientation moduli (ϕ, α, β) of the zigzag to seek the minimum of the NR chemical potential (5.20) for any given combination of \mathbf{F} and M_3/M_4 .

We begin by changing variables from ϵ to $\xi = \epsilon/D$, combining eqs. (5.20) and (5.17), and rewriting the result as

$$\hat{\mu} = \mathbf{F} \times D + N_c \lambda M M_2^2 \times D^2 \xi^2 + \frac{\pi^2 N_c}{20 \lambda M} \times \frac{\mathcal{F}_m(\xi; M_3/M_4)}{D^2} \quad (5.21)$$

where

$$\mathcal{F}_m(\xi; M_3/M_4) = \min_{\alpha, \beta, \phi} \mathcal{F}(\alpha, \beta, \phi; (\xi = \epsilon/D), M_3/M_4) \quad (5.22)$$

for \mathcal{F} from eq. (5.18). Note that although we know \mathcal{F} as an analytic function of all its arguments, and we know how to minimize it analytically with respect to the α and β angles (cf. appendix A), the minimization with respect to the ϕ angle has to be done numerically, so we do not have an explicit formula for the $\mathcal{F}_m(\xi; M_3/M_4)$. Nevertheless, it is fairly easy to calculate numerically both the \mathcal{F}_m function itself and its derivative $\partial \mathcal{F}_m / \partial \xi$.

Before we seek the global minimum of the chemical potential (5.21), let's find all the local extrema. Spelling out the extremality conditions

$$\left(\frac{\partial \mathcal{F}_m}{\partial D} \right)_\xi = 0, \quad \left(\frac{\partial \mathcal{F}_m}{\partial \xi} \right)_D = 0, \quad (5.23)$$

we obtain

$$\frac{\pi^2 N_c}{20 \lambda M} \times \frac{2 \mathcal{F}_m(\xi; M_3/M_4)}{D^3} - N_c \lambda M M_2^2 \times 2 D \xi^2 = \mathbf{F} \quad (5.24)$$

and

$$\frac{\pi^2 N_c}{20\lambda M} \times \frac{1}{D^2} \times \frac{\partial \mathcal{F}_m}{\partial \xi} + N_c \lambda M M_2^2 \times 2D^2 \xi = 0. \quad (5.25)$$

The last equation here has two branches of solutions: the *zigzag branch*

$$\xi > 0, \quad D^4 = \frac{\pi^2}{20(\lambda M M_2)^2} \times \frac{-1}{2\xi} \frac{\partial \mathcal{F}_m}{\partial \xi}, \quad (5.26)$$

and the *straight-line branch*

$$\xi = 0, \quad \text{any } D > 0 \quad (5.27)$$

which follows from $\partial \mathcal{F}_m / \partial \xi = 0$ for $\xi = 0$. Plugging these two solution branches into the compression force equation (5.24), we arrive at

$$\mathbf{F}(D) = N_c \sqrt{\frac{\pi}{\sqrt{20}} \lambda M M_2^3} \times \frac{2\mathcal{F}_m(\xi=0) = 4}{\left[D \times \sqrt{\frac{\sqrt{20}}{\pi} \lambda M M_2} \right]^3} \quad \text{for the straight-line branch,} \quad (5.28a)$$

$$\mathbf{F}(\xi) = N_c \sqrt{\frac{\pi}{\sqrt{20}} \lambda M M_2^3} \times \frac{2\mathcal{F}_m + \xi \frac{\partial \mathcal{F}_m}{\partial \xi}}{\left[\frac{-1}{2\xi} \frac{\partial \mathcal{F}_m}{\partial \xi} \right]^{3/4}} \quad \text{for the zigzag branch.} \quad (5.28b)$$

Between the two branches, for any particular value of the compression force \mathbf{F} we may have 1, 2, or more solutions for the lattice moduli (D, ξ) . Only one of these solutions is the global minimum of the chemical potential (5.20) — the rest are merely local extrema — and the simplest way to identify that global minimum is to evaluate

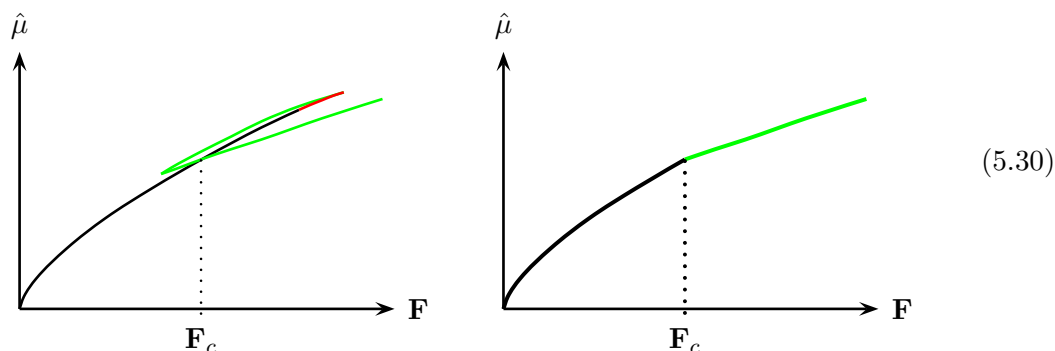
$$\hat{\mu}(D) = N_c M_2 \frac{\pi}{\sqrt{20}} \times \frac{6}{\left[D \times \sqrt{\frac{\sqrt{20}}{\pi} \lambda M M_2} \right]^2} \quad \text{for the straight-line branch,} \quad (5.29a)$$

$$\hat{\mu}(\xi) = N_c M_2 \frac{\pi}{\sqrt{20}} \times \frac{3\mathcal{F}_m + \frac{\xi}{2} \frac{\partial \mathcal{F}_m}{\partial \xi}}{\left[\frac{-1}{2\xi} \frac{\partial \mathcal{F}_m}{\partial \xi} \right]^{1/2}} \quad \text{for the zigzag branch.} \quad (5.29b)$$

for all the solutions of eqs. (5.28) and select the solution with the lowest $\hat{\mu}$.

When we vary the compression force \mathbf{F} or the M_3/M_4 ratio of the background, sometimes the global minimum ‘jumps’ from one solution to another; physically, this corresponds to a phase transition between different lattice geometries. To see how this works, let’s pick an M_3/M_4 ratio — say, $M_3/M_4 = 0.5$ — and plot the chemical potential $\hat{\mu}$ as a function of the compression force \mathbf{F} . for all the solutions of eqs. (5.28). In practice, this means combining the parametric plots $(\mathbf{F}(D), \hat{\mu}(D))$ for the straight-line branch (the black curve) and $(\mathbf{F}(\xi), \hat{\mu}(\xi))$ for the zigzag branch (the red and green curves, according to the zigzag’s

orientation phase), thus



The left plot here shows all the solutions while the right plot shows only the global minimum of the chemical potential $\hat{\mu}$ for any compression force \mathbf{F} . We see that at $\mathbf{F} = \mathbf{F}_c$ the global minimum switches from one solution branch to another; physically, this corresponds to a first-order phase transition from a straight line to a zigzag.

For other values of the M_3/M_4 ratio, we get more complicated $(\mathbf{F}, \hat{\mu})$ plots than (5.30). For example, for $M_3/M_4 = 0.73$ the zigzag branch goes back and forth several times — although this is very hard to see graphically without zooming on very small portions of the curve — so that for some values of the force \mathbf{F} we get up to six different solutions of eqs. (5.28). Consequently, when we vary \mathbf{F} there is a *sequence of three phase transitions*: first, a second-order transition from a straight line to a zigzag with the antiferromagnetic order of orientations. Second, a first order transition in which the zigzag amplitude drastically increases, the lattice spacing shrinks about 20%, and the orientation pattern changes from AF to the non-abelian NA2. Third, another first-order phase transition, this time with a smaller jump of the lattice geometry while the orientation pattern changes from the NA2 to another non-abelian phase NA1.

Altogether, we have found 7 different transition sequences for different M_3/M_4 ratios! But instead of spelling out all these sequences in gory detail, let us simply present in figures 6(a), 6(b), and 7 the phase diagram of all the zigzag and straight-chain phases in three different planes: the compression force \mathbf{F} versus M_3/M_4 , the non-relativistic chemical potential $\hat{\mu}$ versus M_3/M_4 , and the linear instanton density $\rho = 1/D$ versus M_3/M_4 .

Here are a few particularly noteworthy features of these diagrams:

- Since we assume $M_2 \ll M_3$, the straight-chain phase always has the antiferromagnetic order of the instantons' orientations.
- The very first transition from the straight chain to a zigzag could be either first-order or second-order, depending on the M_3/M_4 ratio: for $(M_3/M_4) < 0.725$ the transition is first-order while for $(M_3/M_4) > 0.725$ it's second-order. This difference is due to different orientation phases of the zigzag immediately after the transition: for $(M_3/M_4) > 0.725$ the zigzag has the same antiferromagnetic order as the straight chain, which allows a second-order transition; but for $(M_3/M_4) < 0.725$ the zigzag has a different orientation pattern — the non-abelian NA1 or NA2 — so the transition is first-order.

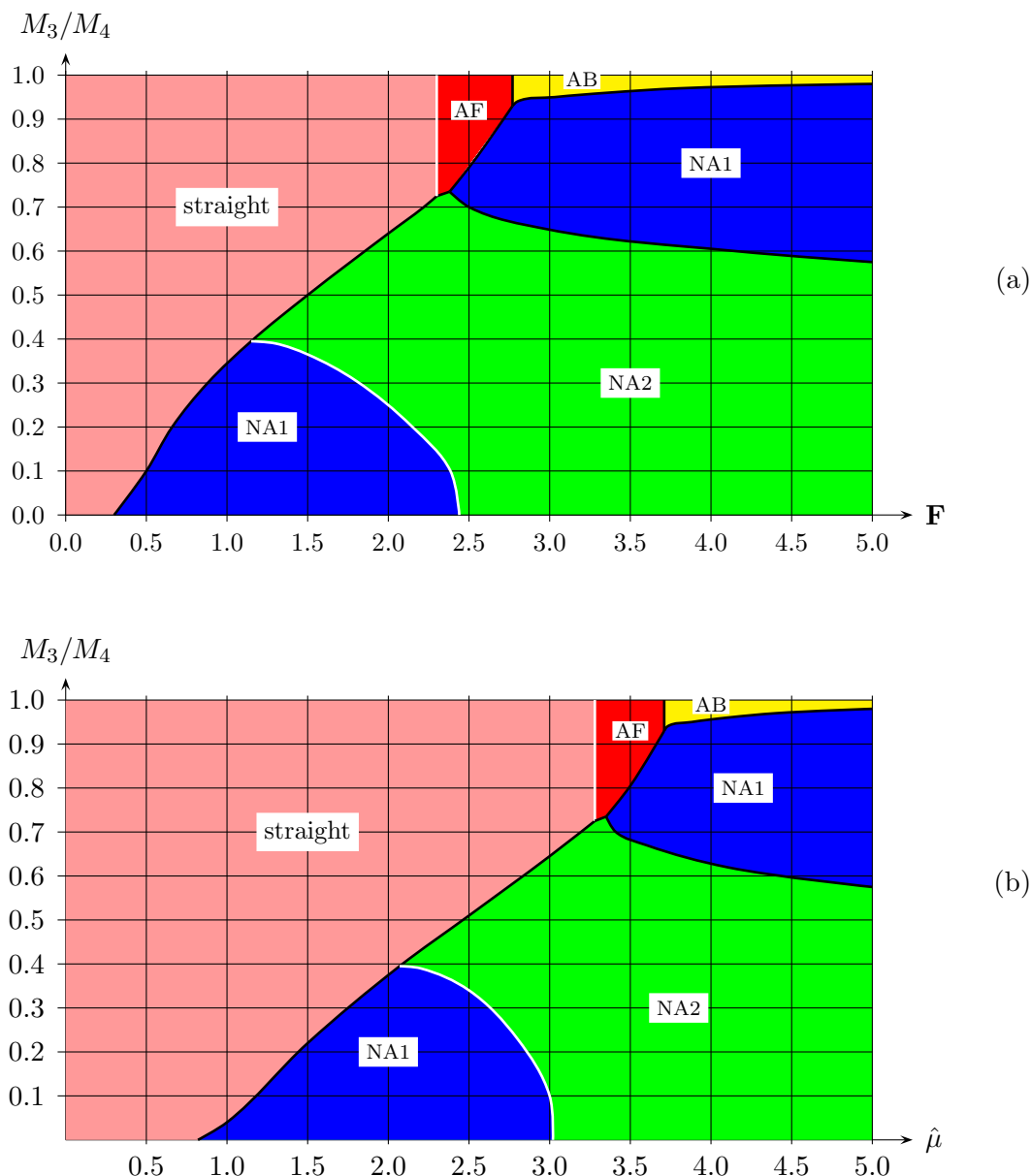


Figure 6. Phase diagrams of the zigzag and straight-chain phases in the compression v. M_3/M_3 plane (top diagram (a), F in units of $N_c\sqrt{\lambda M M_2^3}$) and in the chemical potential v. M_3/M_3 plane (bottom diagram (b), the non-relativistic $\hat{\mu}$ in units of $N_c M_2$). The straight-chain phase is colored pink, while 4 other colors — red, yellow, blue, and green — denote zigzag phases with different instanton orientation patterns. The first-order transition between phases are indicated by black lines, the second-order transitions by white lines.

- The non-abelian phases **NA1** and **NA2** of the zigzag cover much larger areas of the phase diagrams 6(a,b) and 7 than the abelian phases **AF** and **AB**. In particular, at larger compression forces F — and hence larger chemical potentials $\hat{\mu}$, larger densities, and larger zigzag amplitudes — the instanton orientations usually prefer the non-abelian patterns. Only the backgrounds with $M_3 \approx M_4$ — such as the model we have analyzed in [1] — favor the abelian orientations.

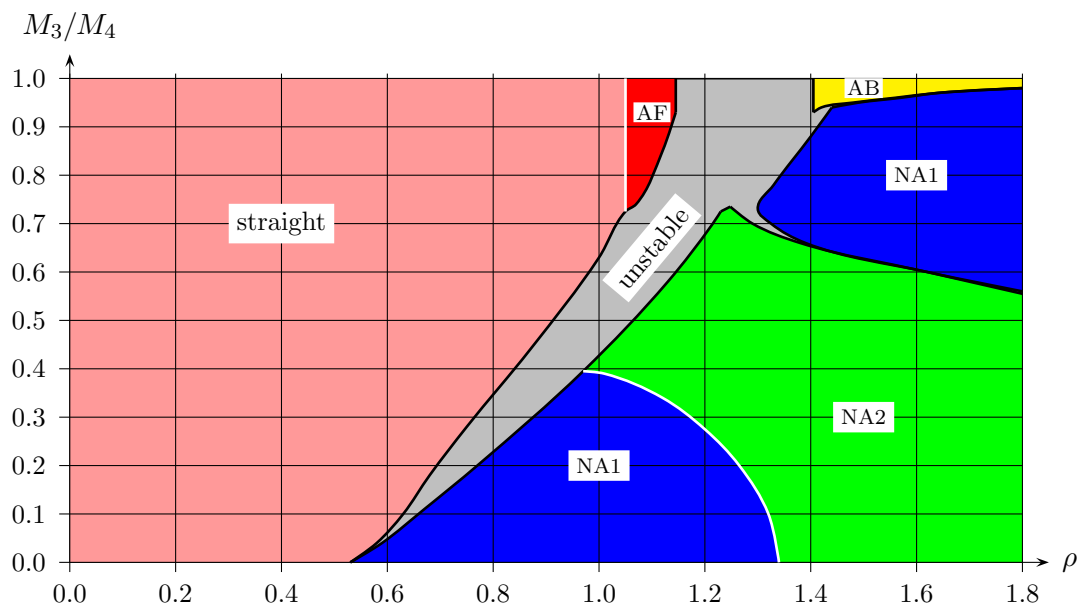


Figure 7. Phase diagram of the zigzag and straight-chain phases in the linear density v . M_3/M_4 plane; the density $\rho = 1/D$ is in units of $\sqrt{\lambda M M_2}$. The stable straight-chain phase is colored pink, while the stable zigzag phases are colored red, yellow, blue, or green according to the instanton orientation pattern. Finally, the gray color denotes densities at which a uniform zigzag or straight chain would be mechanically or thermodynamically unstable.

- Figure 7 has gray areas at which an instanton zigzag with a uniform lattice spacing $D = 1/\rho$ and a uniform amplitude ϵ (or a uniform straight chain for $\epsilon = 0$) would be unstable against instantons' motion along the x^1 axis (the long direction of the zigzag). If we put $N \gg 1$ instantons into a box of fixed length $L = N/\rho$ and let them seek the lowest-energy configuration, they would organize themselves into domains of two different lattices with different lattice spacings and different amplitudes.
- The NA1 phase of the zigzag occupies two separate regions of the phase diagram separated by the region of the other non-abelian phase NA2. The phase transition between the lower-left NA1 region and the NA2 region is second-order, while the transition between the upper-right NA1 region and the NA2 is weakly first-order: the lattice spacing and the zigzag amplitude are discontinuous across the transition, but the discontinuity is very small and hard to see graphically on figure 7 or (5.30)-like plots.
- Likewise, the transition between the upper-right region of the NA1 phase and the abelian AB phases of the zigzag is weakly first-order. On the other hand, the transitions between the antiferromagnetic phases of the straight chain or zigzag and all the other zigzag phases is strongly first-order, with largish discontinuities of the lattice spacing and even larger discontinuity of the zigzag amplitude.

- However, for small M_3/M_4 the discontinuity becomes small; for $M_3/M_4 = 0$ it vanishes altogether and the phase transition between the straight chain and the NA1 phase of the zigzag becomes second-order. This is OK because for $M_3/M_4 = 0$ — or rather for $M_3 = M_2 \ll M_4$ — the instantons' orientations in the straight-chain phase are no longer antiferromagnetic but form the period = 4 Klein-group pattern similar to the NA1 pattern of the zigzag.

The phase diagrams on figures 6(a,b) and 7 completes our analysis of the instanton zigzag. But the zigzag is only the first step in the transition sequence from a 1D to a 2D instanton lattice. The subsequent steps — from a zigzag to a 3-layer lattice, to 4-layer lattice, etc. — will be explored in our next paper [58].

6 Summary and outlook

In [1] we have investigated the crystalline structure of the holographic nuclear matter. The analysis done in that paper revealed the “popcorn transition” of a one-dimensional infinite chain of nucleons into a zigzag structure. We have also found out that the phases between adjacent nucleons are abelian. In this paper we continued the exploration of 1D instanton systems and their popcorn transition to the zigzag, but by generalizing the background to allow for $M_3 \neq M_4$, we found a much richer phase structure, including two distinct non-abelian phases.

In section 3 we proved that for low-density multi-instanton systems — in which the distances between the instantons are much larger than their radii — the interactions between the baryons are dominated by the two-body forces, while the 3-body forces, etc., are suppressed by powers of (radius/distance)². The two-body-force approximation makes for much easier calculations, so in sections 4, 5 we could put the instanton chains and zigzags into richer backgrounds than we had in our previous paper. In particular, we could use a less symmetric external potential for the baryons, by making the 5D gauge coupling vary as

$$\frac{8\pi^2}{g_5^2(x)} = N_c \lambda M \times \left(1 + M_2^2 \times x_2^2 + M_3^2 \times x_3^2 + M_4^2 \times x_4^2 + O(M^4 x_{2,3,4}^4) \right) \quad (1.2)$$

with three independent curvature parameters M_2, M_3, M_4 .

For the almost-holographic background $M_2, M_3 \ll M_4$, we found that the 1D lattice of instantons has a highly degenerate family of ground states with many different patterns of instantons' orientations. Among them are periodic patterns in which instanton orientations span finite subgroups of the $SU(2)/\mathbb{Z}_2$: the anti-ferromagnetic chains spanning the \mathbb{Z}_2 subgroup (two alternating orientations related by a 180° twist), period = 4 chains spanning the Klein group $\mathbb{Z}_2 \times \mathbb{Z}_2$, and period = $2k$ chains spanning the prismatic groups $\mathbb{Z}_k \times \mathbb{Z}_2$ and the dihedral groups D_{2k} . There are also many link-periodic patterns in which the relative orientations $y_n^\dagger y_{n+1}$ of neighboring instantons are periodic but the orientations y_n themselves are not periodic. But the vast majority of the degenerate ground states are not periodic at all. Raising the M_2 and M_3 parameters lifts the degeneracy. For $M_2 \ll M_3 \sim M_4$, the ground state is the anti-ferromagnetic chain, while for $M_2 \sim M_3 \sim M_4$

the ground state is periodic with a longer period or link-periodic, depending on the M_2/M_3 ratio. Specifically, for $M_2 = M_3$ the instanton's orientation in the ground state span the Klein group, for other rational M_2^2/M_3^2 they span a dihedral group, while the irrational M_2^2/M_3^2 favor the link-periodic patterns.

For the zigzag-shaped chains of instantons we need $M_2 \ll M_3 \sim M_4$ to make sure the transition from a straight chain to a zigzag happens for lattice spacing \gg instanton radius, but we may vary the M_3/M_4 ratio. Consequently, we found 5 distinct phases shown in figures 6 and 7: the anti-ferromagnetic straight chain, the anti-ferromagnetic zigzag, the abelian link-periodic zigzag (period = 1), and two different non-abelian link-periodic phases (period = 2). Figures 6 and 7 also show a rather complicated phase structure: as we increase the compression force of the 1D lattice — and hence increase the instanton density and the chemical potential, — the zigzag amplitude and the orientation pattern go through a sequence of 1, 2, or 3 phase transitions, but the number of transition and their thermodynamic orders depend on the M_3/M_4 ratio. In particular, the very first transition from a straight chain to a zigzag is first-order for $M_3 < 0.725M_4$ but second order for $M_3 > 0.725M_4$. The subsequent transition(s) may also be first-order or second-order, depending on the M_3/M_4 .

This paper is part of a program focused on exploring the holographic nuclear matter, especially its low-temperature high-pressure phases. Thus far we have worked out the 1D instanton lattices and their zigzag deformations as a toy model for the transitions between 3D and 4D lattices. But many interesting questions remain open, and we would like to address them in our future papers; here is a short list:

- The very next step in our program is to follow up the “popcorn transitions” from a 1D instanton chain into the second dimension beyond the zigzag phases through multiple lines of instantons towards a thick 2D lattice. We should also check that that the very first popcorn transition indeed goes from the straight chain to the zigzag rather than to some other configuration of the instantons.
- After that, we should set $M_2 = 0$ and study the infinite 2D lattices and their popcorn transitions into the third dimension for $M_3 \ll M_4$. The techniques developed in this paper — using the two-body-forces approximation for $a \ll D$ — should work all the way up to infinite 3D lattices for $M_2 = M_3 = 0$. Alas, the ultimate popcorn transition from the 3D instanton crystal into the fourth space dimension probably happens for $D \sim a$, so the two-body-forces approximation would not be accurate. But we ought to try it anyhow, just to get a qualitative picture of the transition.
- Our program (starting with [1]) is based on the conjecture that the popcorn transition from a 3D instanton crystal to a 4D crystal is holographically dual to the transition of the large- N_c nuclear matter from the baryonic phase to the “quarkyonic phase” [19] (Fermi liquid of quarks, with baryon-like excitations near the Fermi surface). Thus far, this conjecture was justified by qualitative arguments only, and we would like to make a quantitative argument by working out the details and the implication of the transition.

- In the skyrmion model of large- N_c nuclear matter, Kugler and Shtrikman [25, 26] showed that at low pressures, the lowest-energy configuration of is the FCC lattice of skyrmions with 4 different isospin orientations, while the high-pressure configuration is the simple cubic lattice of half-skyrmions. Later, Manton and Sutcliffe [18] described these lattices in terms of instantons on a T^4 torus. We would like to re-interpret their results in terms of instanton lattices to see where the Klebanov-Kugler-Shtrikman transition point (between whole-skyrmion and half-skyrmion lattices) corresponds to the popcorn transition from a 3D to a 4D lattice. Also, we would like to see how the chiral symmetry restoration in the high-density phase works in terms of the instanton lattices.
- Finally, we would like to see if any of our 1D, 2D, 3D lattices — or their generalizations — appear in condensed matter rather than in the holographic nuclear physics. We imagine a crystalline lattice of some high-spin atoms / ions / whatever whose orientations may form all kinds of complicated patterns — perhaps non-abelian patterns — depending on external parameters such as temperature, pressure, or doping. Hopefully, the methods developed in this article would be useful for studying such systems. In this case, we would end up answering questions that have yet not been asked.

Acknowledgments

We thank Dmitry Melnikov for many fruitful conversations. This paper is based on research supported by the US National Science Foundation (grant PHY-0969020) [VK] and by the Israel Science Foundation (grant 1665/10).

A Energy of the instanton zigzag

In this appendix we calculate the net energy of a zigzag-shaped chain (5.2) of small instantons whose orientations follow the ansatz (5.15) from section 5,

$$y_n = \exp(in\frac{\phi}{2}\tau_2) \times \exp\left(i\left(\frac{\alpha}{2} + (-1)^n\frac{\beta}{2}\right)\tau_1\right). \quad (5.15),$$

and then we minimize that energy with respect to the orientation moduli α , β , and ϕ .

The energy of a general 2D instanton configuration in a background with $M_2 \ll M_3 \sim M_4$ is spelled out in eq. (5.9):

$$\mathcal{E} = N_c \lambda M M_2^2 \times \sum_n (X_n^2)^2 + \frac{N_c}{5\lambda M} \sum_{m \neq n} \frac{Q_z(m, n)}{|X_m - X_n|^2} \quad (A.1)$$

for

$$\begin{aligned} Q_z(m, n) = & \frac{1}{2} + \text{tr}^2(y_m^\dagger y_n) + C_3 \sum_{a=1,2} \text{tr}^2(y_m^\dagger y_n (-i\tau^a)) \\ & + (1 - 2C_3) \text{tr}^2(y_m^\dagger y_n (-i\vec{\tau} \cdot \vec{N}_{mn})), \end{aligned} \quad (A.2)$$

where $C_3 = M_3^2/(M_4^2 + M_3^2)$, $0 \leq C_3 \leq \frac{1}{2}$, and $N_{mn}^\mu = (X_m^\mu - X_n^\mu)/|X_m - X_n|$. For the zigzag arrangement of the instanton centers,

$$\left. \begin{aligned} |X_m - X_n|^2 &= (n - m)^2 D^2, \\ \vec{\tau} \cdot \vec{N}_{m,n} &= \pm \tau_1 \end{aligned} \right\} \quad \text{for even } n - m, \quad (\text{A.3})$$

but

$$\left. \begin{aligned} |X_m - X_n|^2 &= (n - m)^2 D^2 + 4\epsilon^2, \\ \vec{\tau} \cdot \vec{N}_{m,n} &= \frac{(n - m)D\tau_1 + (-1)^n(2\epsilon)\tau_2}{\sqrt{(n - m)^2 D^2 + 4\epsilon^2}} \end{aligned} \right\} \quad \text{for odd } n - m. \quad (\text{A.4})$$

Our first task is to calculate the orientation-dependent force coefficients $Q_z(m, n)$ for the instantons oriented according to the ansatz (5.15). For even $\ell = n - m$, we have

$$\begin{aligned} y_m^\dagger y_n &= \exp\left(-i\frac{\alpha \pm \beta}{2}\tau_1\right) \times \exp\left(i\ell\frac{\phi}{2}\tau_2\right) \times \exp\left(+i\frac{\alpha \pm \beta}{2}\tau_1\right) \\ &= \cos\frac{\ell\phi}{2} + \sin\frac{\ell\phi}{2}\cos(\alpha \pm \beta) \times (i\tau_2) + \sin\frac{\ell\phi}{2}\sin(\alpha \pm \beta) \times (i\tau_3) \end{aligned} \quad (\text{A.5})$$

where $\pm = (-1)^n = (-1)^m$, and consequently

$$Q_z(m, n) = \frac{1}{2} + 4\cos^2\frac{\ell\phi}{2} + C_3 \times 4\sin^2\frac{\ell\phi}{2}\cos^2(\alpha \pm \beta). \quad (\text{A.6})$$

Averaging between even and odd n for fixed distance $\ell = n - m$, we obtain

$$\begin{aligned} \langle \cos^2(\alpha \pm \beta) \rangle &\equiv \frac{1}{2} \times \cos^2(\alpha + \beta) + \frac{1}{2} \times \cos^2(\alpha - \beta) \\ &= \cos^2\alpha \cos^2\beta + \sin^2\alpha \sin^2\beta \end{aligned} \quad (\text{A.7})$$

and hence

$$\begin{aligned} \langle Q_z(\ell = n - m) \rangle &= \frac{1}{2} + 4\cos^2\frac{\ell\phi}{2} + 4C_3 \times (\cos^2\alpha \cos^2\beta + \sin^2\alpha \sin^2\beta) \times \sin^2\frac{\ell\phi}{2} \\ &= \frac{9}{2} + 2\left(-1 + C_3(\cos^2\alpha \cos^2\beta + \sin^2\alpha \sin^2\beta)\right) \times (1 - \cos(\ell\phi)). \end{aligned} \quad (\text{A.8})$$

Therefore, the average net energy of an instanton $\#m$ due to 2-body repulsive forces from the other instantons $\#n$ at even distances $\ell = n - m$ comes up to

$$\mathcal{E}_{\text{even}} = \frac{N_c}{5\lambda M} \sum_{\text{even } \ell \neq 0} \frac{\langle Q_z(\ell) \rangle}{(\ell D)^2} \quad (\text{A.9})$$

$$\begin{aligned} &= \frac{N_c}{5\lambda M D^2} \times \left(\frac{9}{2} \times \left(\sum_{\text{even } \ell \neq 0} \frac{1}{\ell^2} = \frac{\pi^2}{12} \right) \right. \\ &\quad \left. + 2\left(-1 + C_3(\cos^2\alpha \cos^2\beta + \sin^2\alpha \sin^2\beta)\right) \times \left(\sum_{\text{even } \ell \neq 0} \frac{1 - \cos(\ell\phi)}{\ell^2} = \frac{\pi^2}{8} \Sigma_0(\phi) \right) \right) \\ &= \frac{\pi^2 N_c}{20\lambda M D^2} \left(\frac{3}{2} + \left(-1 + C_3(\cos^2\alpha \cos^2\beta + \sin^2\alpha \sin^2\beta)\right) \times \Sigma_0(\phi) \right) \end{aligned} \quad (\text{A.10})$$

where

$$\Sigma_0(\phi) = \frac{4\phi(\pi - \phi)}{\pi^2} \quad (\text{A.11})$$

cf. eq. (5.19a).

For instantons at odd distances from each other we have more complicated formulae. For odd $\ell = n - m$ we have

$$\begin{aligned} y_m^\dagger y_n &= \exp\left(\frac{-i\alpha \pm i\beta}{2} \tau_1\right) \times \exp\left(i\ell \frac{\phi}{2} \tau_2\right) \times \exp\left(\frac{+i\alpha \pm i\beta}{2} \tau_1\right) \\ &= \cos \frac{\ell\phi}{2} \times \exp(\pm i\beta \tau_1) + \sin \frac{\ell\phi}{2} \times (i\tau_2) \times \exp(i\alpha \tau_1) \\ &= \cos \frac{\ell\phi}{2} \cos \beta \pm \cos \frac{\ell\phi}{2} \sin \beta \times (i\tau_1) + \sin \frac{\ell\phi}{2} \cos \alpha \times (i\tau_2) + \sin \frac{\ell\phi}{2} \sin \alpha \times (i\tau_3) \end{aligned} \quad (\text{A.12})$$

where $\pm = (-1)^n = -(-1)^m$. Consequently,

$$\text{tr}(y_m^\dagger y_n (-i\vec{\tau} \cdot \vec{N}_{mn})) = \frac{\pm 2}{\sqrt{(\ell D)^2 + 4\epsilon^2}} \times \left(\cos \frac{\ell\phi}{2} \sin \beta \times \ell D - \sin \frac{\ell\phi}{2} \cos \alpha \times 2\epsilon \right) \quad (\text{A.13})$$

and hence

$$\begin{aligned} Q_z(m, n) &= \frac{1}{2} + 4 \cos^2 \frac{\ell\phi}{2} \cos^2 \beta + C_3 \left(4 \cos^2 \frac{\ell\phi}{2} \sin^2 \beta + 4 \sin^2 \frac{\ell\phi}{2} \cos^2 \alpha \right) \\ &\quad + \frac{4(1 - 2C_3)}{\ell^2 D^2 + 4\epsilon^2} \times \left(\cos \frac{\ell\phi}{2} \sin \beta \times \ell D - 2 \sin \frac{\ell\phi}{2} \cos \alpha \times 2\epsilon \right)^2. \end{aligned} \quad (\text{A.14})$$

Note that this $Q_z(m, n)$ depends only on $\ell = n - m$, so there is no need to average it over m . Instead, it is convenient to expand the right hand side of this formula and to re-group the terms according to their ℓ dependence, thus

$$\begin{aligned} Q_z(\ell) &= \left(\frac{5}{2} - \sin^2 \beta + \cos^2 \alpha \right) + (2 - \sin^2 \beta - \cos^2 \alpha) \times \cos(\ell\phi) \\ &\quad + (1 - 2C_3)(\sin^2 \beta - \cos^2 \alpha) \times \frac{\ell^2 D^2 - 4\epsilon^2}{\ell^2 D^2 + 4\epsilon^2} \\ &\quad + (1 - 2C_3)(\sin^2 \beta + \cos^2 \alpha) \times \frac{\ell^2 D^2 - 4\epsilon^2}{\ell^2 D^2 + 4\epsilon^2} \cos(\ell\phi) \\ &\quad - 2(1 - 2C_3) \cos \alpha \sin \beta \times \frac{4\ell D \epsilon}{\ell^2 D^2 + 4\epsilon^2} \sin(\ell\phi). \end{aligned} \quad (\text{A.15})$$

Therefore, the net energy per instanton due to repulsion from its odd-distance neighbors

$$\mathcal{E}_{\text{odd}} = \frac{N_c}{5\lambda M} \sum_{\text{odd } \ell} \frac{Q_z(\ell)}{\ell^2 D^2 + 4\epsilon^2} \quad (\text{A.16})$$

is a linear combination of five series,

$$\mathcal{E}_{\text{odd}} = \frac{\pi^2 N_c}{20\lambda M D^2} \left(\begin{aligned} &\left(\frac{5}{2} - \sin^2 \beta + \cos^2 \alpha \right) \times \Sigma_1 + (2 - \sin^2 \beta - \cos^2 \alpha) \times \Sigma_2 \\ &\quad + (1 - 2C_3)(\sin^2 \beta - \cos^2 \alpha) \times \Sigma_3 \\ &\quad + (1 - 2C_3)(\sin^2 \beta + \cos^2 \alpha) \times \Sigma_4 \\ &\quad - 2(1 - 2C_3) \cos \alpha \sin \beta \times \Sigma_5 \end{aligned} \right) \quad (\text{A.17})$$

where

$$\begin{aligned}
 \Sigma_1 &\equiv \frac{4D^2}{\pi^2} \sum_{\text{odd } \ell} \frac{1}{\ell^2 D^2 + 4\epsilon^2} = \frac{\tanh(\pi\epsilon/D)}{(\pi\epsilon/D)}, \\
 \Sigma_2 &\equiv \frac{4D^2}{\pi^2} \sum_{\text{odd } \ell} \frac{\cos \ell\phi}{\ell^2 D^2 + 4\epsilon^2} = \frac{\sinh((\pi - 2\phi)\epsilon/D)}{(\pi\epsilon/D) \cosh(\pi\epsilon/D)}, \\
 \Sigma_3 &\equiv \frac{4D^2}{\pi^2} \sum_{\text{odd } \ell} \frac{\ell^2 D^2 - 4\epsilon^2}{[\ell^2 D^2 + 4\epsilon^2]^2} = \frac{1}{\cosh^2(\pi\epsilon/D)}, \\
 \Sigma_4 &\equiv \frac{4D^2}{\pi^2} \sum_{\text{odd } \ell} \frac{(\ell^2 D^2 - 4\epsilon^2) \times \cos \ell\phi}{[\ell^2 D^2 + 4\epsilon^2]^2} = \frac{\cosh(2\phi\epsilon/D)}{\cosh^2(\pi\epsilon/D)} - \frac{2\phi}{\pi} \times \frac{\cosh((\pi - 2\phi)\epsilon/D)}{\cosh(\pi\epsilon/D)}, \\
 \Sigma_5 &\equiv \frac{4D^2}{\pi^2} \sum_{\text{odd } \ell} \frac{4\ell D\epsilon \times \sin \ell\phi}{[\ell^2 D^2 + 4\epsilon^2]^2} = \frac{\sinh(2\phi\epsilon/D)}{\cosh^2(\pi\epsilon/D)} + \frac{2\phi}{\pi} \times \frac{\sinh((\pi - 2\phi)\epsilon/D)}{\cosh(\pi\epsilon/D)},
 \end{aligned} \tag{A.18}$$

cf. eqs. (5.19b)–(5.19f).

Altogether, combining the interaction energies between even-distance and odd-distance neighbors we obtain the net average energy of an instanton in a zigzag,

$$\mathcal{E}_{\text{per instanton}}^{\text{interaction}} = \mathcal{E}_{\text{even}} + \mathcal{E}_{\text{odd}} = \frac{\pi^2 N_c}{20\lambda M D^2} \times \mathcal{F}(\phi, \alpha, \beta) \tag{A.19}$$

for

$$\begin{aligned}
 \mathcal{F} = & \frac{3}{2} + \left(-1 + C_3(\cos^2 \alpha \cos^2 \beta + \sin^2 \alpha \sin^2 \beta) \right) \times \Sigma_0(\phi) \\
 & + \left(\frac{5}{2} + \cos^2 \alpha - \sin^2 \beta \right) \times \Sigma_1(\epsilon/D) \\
 & + \left(2 - \cos^2 \alpha - \sin^2 \beta \right) \times \Sigma_2(\phi, \epsilon/D) \\
 & + (1 - 2C_3)(\sin^2 \beta - \cos^2 \alpha) \times \Sigma_3(\epsilon/D) \\
 & + (1 - 2C_3)(\sin^2 \beta + \cos^2 \alpha) \times \Sigma_4(\phi, \epsilon/D) \\
 & - 2(1 - 2C_3) \cos \alpha \sin \beta \times \Sigma_5(\phi, \epsilon/D).
 \end{aligned} \tag{5.18}$$

Or rather, this is the net zigzag's energy (per instanton) due to 2-body interactions, but there is also a 1-body potential energy

$$\mathcal{E}_{\text{pot}} = N_c \lambda M M_2^2 \times \epsilon^2 \tag{A.20}$$

due to instanton centers being at $X_n^2 = \pm\epsilon \neq 0$. Adding this potential energy to eq. (A.19) we arrive at

$$\mathcal{E}_{\text{per instanton}}^{\text{zigzag}} = \frac{\pi^2 N_c}{20\lambda M D^2} \times \mathcal{F}(\phi, \alpha, \beta) + N_c \lambda M M_2^2 \times \epsilon^2 \tag{A.21}$$

and hence eq. (5.17). *Quod erat demonstrandum.*

★ ★ ★

In the second part of this appendix we minimize the zigzag's energy with respect to the orientation moduli ϕ, α, β . Specifically, we are going to minimize $\mathcal{F}(\phi, \alpha, \beta)$ in two stages:

at first we hold ϕ fixed and minimize WRT α and β , and then we allow ϕ to vary and seek the overall minimum.

For the first stage it is convenient to change variables

$$\alpha \rightarrow a = \cos \alpha \quad \text{and} \quad \beta \rightarrow b = \sin \beta, \quad -1 \leq a, b \leq +1. \quad (\text{A.22})$$

Then in terms of (ϕ, a, b) the \mathcal{F} function becomes a polynomial in a and b with ϕ -dependent coefficients,

$$\mathcal{F}(\phi, a, b) = A(\phi) + B(\phi) \times a^2 + C(\phi) \times b^2 - 2J(\phi) \times ab - K(\phi) \times a^2 b^2 \quad (\text{A.23})$$

where

$$\begin{aligned} A(\phi) &= \frac{3}{2} - \Sigma_0(\phi) + \frac{5}{2}\Sigma_1 + 2\Sigma_2(\phi), \\ B(\phi) &= C_3\Sigma_0(\phi) + \Sigma_1 - \Sigma_2(\phi) - (1 - 2C_3)\Sigma_3 + (1 - 2C_3)\Sigma_4(\phi), \\ C(\phi) &= C_3\Sigma_0(\phi) - \Sigma_1 - \Sigma_2(\phi) + (1 - 2C_3)\Sigma_3 + (1 - 2C_3)\Sigma_4(\phi), \\ J(\phi) &= (1 - 2C_3)\Sigma_5(\phi), \\ K(\phi) &= 2C_3\Sigma_0(\phi). \end{aligned} \quad (\text{A.24})$$

Minimizing the polynomial (A.23) WRT a and b works differently for $\phi = 0$ or $\phi = \pi$ and for generic ϕ 's, so let us start with the special cases. For $\phi = \pi$ we have $\Sigma_0 = \Sigma_5 = 0$ while $\Sigma_2 = -\Sigma_1$ and $\Sigma_4 = -\Sigma_3$, hence

$$K = J = C = 0 \quad \text{while} \quad B > 0. \quad (\text{A.25})$$

Consequently, \mathcal{F} becomes independent of b , but that's OK because the β angle does not affect the instantons' relative orientations $y_m^\dagger y_n$ for $\phi = \pi$, cf. eqs. (A.5) and (A.12). In other words, β or b becomes an irrelevant variable at $\phi = \pi$ — just like the geographic longitude becomes an irrelevant coordinate at latitudes $\pm 90^\circ$. As to the a variable, $B > 0$ means that the minimum is at $a = 0$, i.e., $\alpha = \pm \frac{\pi}{2}$. Altogether, we have

$$y_n^\dagger y_{n+1} = \pm i\tau_3 \quad (\text{A.26})$$

which means the anti-ferromagnetic phase **AF** of the instanton orientations. The energy of this phase corresponds to

$$\mathcal{F}(\text{AF}) = A(\phi = \pi) = \frac{3}{2} + \frac{1}{2}\Sigma_1. \quad (\text{A.27})$$

Similarly, for $\phi = 0$ we have $\Sigma_0 = \Sigma_5 = 0$ while $\Sigma_2 = +\Sigma_1$ and $\Sigma_4 = +\Sigma_3$, hence

$$K = J = B = 0 \quad \text{while} \quad C < 0. \quad (\text{A.28})$$

This time, \mathcal{F} becomes independent of a , but that's OK since the α angle becomes irrelevant to the instanton's relative orientations $y_m^\dagger y_n$ for $\phi = 0$, cf. eqs. (A.5) and (A.12). As to the b variable, $C < 0$ calls for maximal $b^2 = \sin^2 \beta$, thus $b = \pm 1$ and $\beta = \pm \frac{\pi}{2}$. Altogether we obtain

$$y_n^\dagger y_{n+1} = \pm i\tau_1 \quad (\text{A.29})$$

and hence a new anti-ferromagnetic phase AF' . However, this new anti-ferromagnetic phase AF' has a higher energy than the old anti-ferromagnetic phase AF —

$$\mathcal{F}(\text{AF}') = A(\phi = 0) + C(\phi = 0) = \frac{3}{2} + \frac{5}{2}\Sigma_1 + 2(1 - 2C_3)\Sigma_3 > \mathcal{F}(\text{AF}) \quad (\text{A.30})$$

— so it is not going to survive the minimization of energy WRT ϕ .

Now consider generic values of ϕ , $0 < \phi < \pi$. This time we have

$$K > 0, \quad J \geq 0, \quad B - C > 0, \quad (\text{A.31})$$

although the signs of B and C themselves depend on ϕ and the parameters ϵ/D and M_3/M_4 . Thanks to $K > 0$, the polynomial (A.23) cannot have any minima at generic values of a and b . More precisely, a, b are limited to the square $-1 \leq a, b \leq +1$, and \mathcal{F} cannot have any local minima in the interior of this square (i.e. for $-1 < a, b < +1$) except maybe at the center $a = b = 0$.

To see this, consider any straight line through the center of the square. Along such a line the ratio a/b is fixed, so \mathcal{F} becomes a quadratic polynomial of $\text{radius}^2 = a^2 + b^2$ with a non-positive coefficient of the $(\text{radius}^2)^2$ term. As a function of an un-bounded real variable such a polynomial does not have any minima at all. But since the radius^2 variable is bounded — between zero at the center and the maximum at the edge of the square — \mathcal{F} may have a minimum (or minima) at the center or/and at the edge. Applying this argument to all lines through the center, we find that the minimum or minima of \mathcal{F} may lie at the square's boundaries $a = \pm 1$ or $b = \pm 1$, or at the center $a = b = 0$, but not at any other interior point.

Let us consider such possible minima in more detail:

- The minimum at $a = b = 0$ — i.e. $\alpha = \pm \frac{\pi}{2}$, $\beta = 0$ — leads to

$$y_n^\dagger y_{n+1} = \cos \frac{\phi}{2} \pm \sin \frac{\phi}{2} \times (i\tau_3) \quad (\text{A.32})$$

and hence the abelian AB phase of the instanton orientations. Local stability of this minimum requires $C > 0$ and $BC > J^2$, and its energy corresponds to

$$\mathcal{F}(\text{AB}) = A(\phi). \quad (\text{A.33})$$

- Minimum at the boundary $b = \pm 1$ (but generic $-1 < a < +1$) — i.e. $\beta = \pm \frac{\pi}{2}$ and generic α — leads to

$$y_n^\dagger y_{n+1} = (-1)^n \cos \frac{\phi}{2} \times (\pm i\tau_1) + \sin \frac{\phi}{2} \times \left(\cos \alpha \times (i\tau_2) + \sin \alpha \times (i\tau_3) \right) \quad (\text{A.34})$$

which corresponds to the non-abelian orientation phase NA2 . The specific location of such a minimum is

$$b = \pm 1, \quad a = \pm \frac{J}{B - K} \quad (\text{A.35})$$

and its local stability at $|a| < 1$ requires

$$B - K > J \quad \text{but} \quad \left(\frac{J}{B - K} \right)^2 > \frac{C}{B}. \quad (\text{A.36})$$

The energy of this minimum is

$$\mathcal{F}(\text{NA2}) = A(\phi) + B(\phi) - \frac{J^2(\phi)}{B(\phi) - K(\phi)}. \quad (\text{A.37})$$

- In principle there could be a similar minimum at the other boundary $a = \pm 1$ (but generic b). Specifically, such a minimum would happen at

$$a = \pm 1, \quad b = \pm \frac{J}{C - K} \quad (\text{A.38})$$

and its local stability at $|b| < 1$ would require

$$C - K > J \quad \text{but} \quad \left(\frac{J}{C - K} \right)^2 > \frac{B}{C}. \quad (\text{A.39})$$

However, thanks to $B > C$ these two requirements contradict each other, so this type of minimum does not happen.

- Finally, we may have minima at the corners of the (a, b) square, at $a, b = \pm 1$ i.e., $\alpha = 0$ or π while $\beta = \pm \frac{\pi}{2}$. Such minima lead to

$$y_n^\dagger y_{n+1} = (-1)^n \cos \frac{\phi}{2} \times (\pm i \tau_1) + \sin \frac{\phi}{2} (\pm i \tau_2) \quad (\text{A.40})$$

and hence the non-abelian [NA1](#) phase of the instanton orientations. Local stability of this phase requires $B \geq J + K$ while its energy corresponds to

$$\mathcal{F}(\text{NA1}) = A(\phi) + B(\phi) + C(\phi) - 2J(\phi) - K(\phi). \quad (\text{A.41})$$

Altogether, for each generic $\phi \neq 0, \pi$ we have one or more local energy minima with respect to α and β moduli, and we have analytic formulae [\(A.33\)](#), [\(A.37\)](#), [\(A.41\)](#) for their energies (or rather \mathcal{F} functions) $\mathcal{F}[\text{AB}](\phi)$, $\mathcal{F}[\text{NA2}](\phi)$, $\mathcal{F}[\text{NA1}](\phi)$ as functions of the ϕ modulus. The next step is to minimize these functions with respect to ϕ , but unfortunately the zero-derivative equations

$$\frac{\partial \mathcal{F}[\text{AB}]}{\partial \phi} = 0, \quad \frac{\partial \mathcal{F}[\text{NA2}]}{\partial \phi} = 0, \quad \frac{\partial \mathcal{F}[\text{NA1}]}{\partial \phi} = 0, \quad (\text{A.42})$$

are transcendental and do not have analytic solutions. In principle, we could solve these equations numerically and then seek the global minimum among the solutions, but since we do not seek high precision we have used a simpler numeric method: covered the $0 < \phi < \pi$ interval with a 1800-point grid, evaluated the $\mathcal{F}[\text{AB}]$, $\mathcal{F}[\text{NA2}]$, and $\mathcal{F}[\text{NA1}]$ for all grid points — and also the $\mathcal{F}[\text{AF}]$ for $\phi = \pi$, — and looked for the lowest value among these $3 \times 1800 + 1$ data points.

The phase diagram on figure [5](#) (on page [52](#)) was obtained by repeating this numerical minimization procedure for 100×100 combinations of the ϵ/D and M_3/M_4 parameters.

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